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Region 2 RAC2 Remedial Action Contract

Data Evaluation Report

Pierson's Creek Site, Operable Unit 2

Remedial Investigation/ Feasibility Study Oversight

Newark, New Jersey

June 3, 2019



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Acronyms

ABS absolute difference

CDM Smith CDM Federal Programs Corporation CLP Contract Laboratory Program

CRQL contract required quantitation limit

DQI data quality indicator DQO data quality objective

DMC deuterated monitoring compound

DV data validation

EPA United States Environmental Protection Agency

FS Feasibility Study

Hg mercury

GC/MS gas chromatography/mass spectrometry

ICP-AES inductively coupled plasma-atom emission spectroscopy

ICP-MS inductively coupled plasma-mass spectrometry

ID identification

LCS laboratory control sample

LCSD laboratory control sample duplicate

MS matrix spike

MSD matrix spike duplicate mg/kg milligram per kilogram mg/L milligram per liter

N normal NA not applicable NC not calculable

ng/mL nanogram per milliliter

OU operable unit
PAL project action limit
% D percent difference
% R percent recovery

% RSD percent relative standard deviation

PRP potentially responsible party
PPW Paired Prentice Wilcoxon

QA quality assurance QC quality control

QAPP quality assurance project plan

QL quantitation limit

rev. revision

RI remedial investigation

RL reporting limit

RPD relative percent difference RRF relative response factor RSD relative standard deviation

SO soil

SOP standard operating procedure

SOW statement of work SM standard method



semivolatile organic compound SVOC

Target Analyte List TAL

TB

trip blank Target Compound List Pierson's Creek Site TCL the Site

volatile organic compound VOC

WG groundwater water quality WQ μg/L microgram per liter



Section 1

Introduction

CDM Federal Programs Corporation (CDM Smith) performed oversight of potentially responsible party (PRP) field activities for the Remedial Investigation (RI) / Feasibility Study (FS) at the Pierson's Creek Site (the Site), Operable Unit 2, located in Newark, New Jersey. The overall objective of the field oversight activities was to verify the quality of the PRP's data and ensure that work is performed in compliance with approved plans. This included verification that all field activities were performed in accordance with the PRP's final quality assurance project plan (QAPP) and acceptance and analysis of split samples for comparison with the PRP's data. This report compares CDM Smith's split sample results to the PRP sample results and discusses the differences in the data pairs. In this document, samples are referred to as either PRP samples or CDM Smith split samples for clarity. This report also discusses usability of split samples data accepted by CDM Smith and documents whether those data meet quality objectives and user requirements outlined in the Pierson's Creek Site Final QAPP (CDM Smith 2018).

Sampling activities occurred from December 4, 2018 through December 27, 2018 and February 9, 2019. CDM Smith accepted split samples from soil and groundwater sampling activities at a rate of approximately 10 percent of the PRP's samples by matrix. The split samples were subsequently submitted to the Environmental Protection Agency (EPA) Contract Laboratory Program (CLP) laboratory for analysis. Samples accepted and methods used to analyze the samples are summarized in Table 1-1.



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Section 2

Usability Summary

Split sample data generated during the RI/FS oversight are considered definitive data generated under an EPA approved QAPP, following EPA methods and validated according to EPA Region 2 Guidelines. Data were validated by EPA data validation contractors.

The Data Usability Worksheet discusses field sampling and validation process and summarizes the achievement of the data quality objectives (Table 2-1). Only the final qualified data will be presented in the reports pertaining to the Site. The split sample data that did not meet quality control (QC) criteria were appropriately qualified as estimated (J, J+, J-, JN, or NJ), estimated non-detect (U or UJ), or as rejected (R) during data validation. All data reported herein are usable with the following data validation qualifiers added.

- J: The analyte was positively identified, and the associated numerical value is the approximate concentration of the analyte in the sample.
- J+: The result is an estimated quantity, but the result may be biased high.
- J-: The result is an estimated quantity, but the result may be biased low.
- JN or NJ: The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents its approximate concentration.
- R: The sample results are unusable due to poor data quality; certain criteria were not met.
 The analyte may or may not be present in the sample at the reported concentration.
- U: The analyte was not detected at a level greater than or equal to the level of the sample contract required quantitation limit (CRQL).
- UJ: The analyte was not detected; however, the reported CRQL is approximate and may be inaccurate or imprecise.

Less than one percent (%) of CDM Smith's semivolatile organic compound (SVOC) results were rejected due to surrogate recovery. Two samples, GS-B-112-001-CDM and GS-B-114-001-CDM, had 1,4-dioxane results rejected. Completeness Tables 2-2a and 2-2b show data qualified for environmental samples accepted during the RI/FS oversight split sampling events. Final percentages of valid data are 99.85 % for soil and 100 % for groundwater. Thus, the 90 percent completeness goal for usable data has been met. Table 2-3 presents trip blank results.



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Section 3

Data Quality Indicators

Precision, accuracy, representativeness, comparability, completeness, and sensitivity are used as data quality indicators (DQIs). These DQIs provide a mechanism for evaluating and measuring data quality throughout the project. Achievement of the project's quality objectives was measured from the DQI results assessed against their measurement performance criteria described in Worksheets 12 and 28 of the Final QAPP (CDM Smith 2018). QC parameters evaluated in data review/data validation and corresponding DQIs are discussed in Table 2-1.

As part of the oversight assignment, CDM Smith was tasked to accept split samples for 10 percent of the PRP's soil and groundwater samples; and compare split sample results with the PRP data. A relative percent difference (RPD) was calculated when both CDM Smith split and PRP samples results were greater than five times EPA's CRQL. When either of the results were equal to or less than five times EPA's CRQL, an absolute difference (ABS) was calculated. When both results were not detected, neither RPD nor ABS were calculated.

CDM Smith split samples and PRP samples were evaluated using a criterion of 100 percent RPD for soil samples and 50 percent RPD for groundwater samples. For all other data sets, the ABS should be less than two times EPA's CRQL. The comparison results are provided in Tables 3-1a and 3-1b for soil samples and Table 3-2 for groundwater samples. RPD/ABS exceedances are summarized in Table 2-1; chemicals with RPD/ABS exceedances are summarized below.

- Soil samples:
 - Metals: antimony, arsenic, silver, sodium, thallium, vanadium, and zinc
 - PCBs: Aroclor 1242, Aroclor 1254, and Aroclor 1260
 - SVOCs: 1,1-biphenyl, 1,2,4,5-tetrachlorobenzene, 2-methylnapthalene, acenaphthene, acenaphthylene, anthracene, benzo(a)anthracene, benzo(a)pyrene, benzo(b)fluoranthene, benzo(g,h,i)perylene, benzo(k)fluoranthene, bis(2-ethylhexyl)phthalate, carbazole, chrysene, dibenzo(a,h)anthracene, dibenzofuran, dimethylphthalate, di-n-butylphthalate, di-n-octylphthalate, fluoranthene, fluorene, indeno(1,2,3-cd)pyrene, naphthalene, phenanthrene, phenol, and pyrene
 - VOCs: 1,2,3-trichlorobenzene, 1,2,4-trichlorobenzene, 1,4-dichlorobenzene, 2-butanone, acetone, benzene, carbon disulfide, chlorobenzene, cis-1,2-dichloroethene, cyclohexane, ethylbenzene, isopropylbenzene, m,p-xylene, methyl acetate, methylcyclohexane, methylene chloride, o-xylene, tetrachloroethene, toluene, and trichloroethene
- Groundwater samples:
 - Metals: aluminum, arsenic, iron, lead, mercury, and zinc
 - Pesticides: delta-BHC
 - VOCs: 1,2-dichlorobenzene, cyclohexane, m,p-xylene, methylcyclohexane, and methylene chloride



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Section 4

Data Comparison

Chemicals with RPD/ABS exceedance were further evaluated using a Paired Prentice Wilcoxon (PPW) test. These selected analytes were evaluated only when there were at least four detected sample pairs. The statistical tests were performed to calculate p-values. The p-value indicates whether the difference between the data pairs is statistically significant. A p-value of less than 0.05 indicates a statistically significant difference between the two datasets. PPW tests were conducted to allow inclusion of the left-censored (nondetected) data pairs for selected analytes. The elimination of data pairs containing nondetected values is essentially equivalent to ignoring potentially substantial information contained within these nondetect-containing data pairs and may have led to biased results (Helsel 2005).

4.1 Soil Samples

PPW tests results for soil sample pairs are presented in Table 4-1. Most results were comparable, except the following chemicals:

- Metals: antimony and silver
- PCBs: Aroclor 1254
- SVOCs: benzo(a)anthracene, benzo(a)pyrene, benzo(b)fluoranthene, benzo(g,h,i)perylene, benzo(k)fluoranthene, chrysene, fluoranthene, and indeno(1,2,3-cd)pyrene
- VOCs: o-xylene

The difference in results between PRP and CDM Smith split data pairs was consistent; PRP sample concentrations were generally higher than CDM Smith sample concentrations.

4.2 Groundwater Samples

Only two groundwater split samples were accepted. Thus, PPW tests were not performed on groundwater samples with RPD/ABS exceedances. The difference in results between PRP and CDM Smith data pairs was consistent. PRP sample concentrations were consistently less than CDM Smith sample concentrations for metal and VOCs results. Pesticides (delta-BHC) were only detected in one PRP sample.



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Section 5

References

PRP Documents

Geosyntec Consultants. 2018. Quality Assurance Project Plan, Appendix 1 – Remedial Action Work Plan, Operable Unit 2, Pierson's Creek Superfund Site. Troy Chemical Newark Manufacturing Plan. Newark, New Jersey. March.

CDM Smith Documents

CDM Federal Programs Corporation (CDM Smith). 2018. Final Quality Assurance Project Plan. Pierson's Creek Site, Newark, New Jersey. July 30.

Data Validation Procedures (EPA Region 2 Protocols)

Procedures (SOP) No. HW-33A. Low/Medium Volatile Data Validation. Revision 1. September.
2015. HW-34A. Trace Volatile Data Validation. Revision 1. June.
2016. Hazardous Waste Support Section, SOP No. HW-35A. Semivolatile Data Validation. Revision 1. September.
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2015. Hazardous Waste Support Section, SOP No. HW-37A. Polychlorinated Biphenyl (PCB Aroclor Data Validation. Revision 0. June.
2015. Hazardous Waste Support Section, SOP No. HW-3A. Inductively coupled plasma-ator emission spectroscopy (ICP-AES) Data Validation. Revision 1. July.
2016. Hazardous Waste Support Section, SOP No. HW-3B. Inductively coupled plasma-mas spectroscopy (ICP-MS) Data Validation. Revision 1. September.
2015. Hazardous Waste Support Section, SOP No. HW-3C. Mercury and Cyanide Data Validation. Revision 1. July.

Analytical Methodologies

EPA. 2016. EPA Contract Laboratory Program, Statement of Work for Organic Superfund Methods, Multi-media, Multi-Concentration. SOM02.4. October.

EPA. 2016. EPA Contract Laboratory Program, Statement of Work for Inorganic Superfund Methods, Multi-media, Multi-Concentration. ISM02.4. October.

Data Comparison

Helsel. D.R. 2005. Nondetects and Data Analysis: Statistics for Censored Environmental Data. Wiley-Interscience.



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Tables

Tables



Table 1-1 Summary of Split Samples Pierson's Creek Site, Operable Unit 2 Newark, New Jersey

Location	Sample ID	PRP's Sample ID	Medium	Sample Type	Sample Date	Mercury	Metals	PCBs	Pesticides	SVOCs	VOCs
GS-B-101	GS-B-101-001-CDM	GS-B-101-001-GEOSY	SO	N	12/14/2018	1	22	9		68	
GS-B-101	GS-B-101-004-CDM	GS-B-101-004-GEOSY	SO	N	12/14/2018						51
GS-B-102	GS-B-102-001-CDM	GS-B-102-001-GEOSY	SO	N	12/17/2018	1	22	9		68	
GS-B-102	GS-B-102-002-CDM	GS-B-102-002-GEOSY	SO	N	12/17/2018	1	22	9		68	
GS-B-102	GS-B-102-004-CDM	GS-B-102-004-GEOSY	SO	N	12/17/2018						51
GS-B-106	GS-B-106-004-CDM	GS-B-106-004-GEOSY	SO	N	12/13/2018						51
GS-B-111	GS-B-111-001-CDM	GS-B-111-001-GEOSY	SO	N	12/17/2018	1	22	9		68	
GS-B-111	GS-B-111-004-CDM	GS-B-111-004-GEOSY	SO	N	12/17/2018						51
GS-B-112	GS-B-112-001-CDM	GS-B-112-001-GEOSY	SO	N	12/13/2018	1	22	9		68	
GS-B-112	GS-B-112-004-CDM	GS-B-112-004-GEOSY	SO	N	12/13/2018						51
GS-B-114	GS-B-114-001-CDM	GS-B-114-001-GEOSY	SO	N	2/9/2019	1	22	9		68	51
GS-SS-101	GS-SS-101-001-CDM	GS-SS-101-001-GEOSY	SO	N	12/4/2018	1	22	9		68	
GS-SS-101	GS-SS-101-002-CDM	GS-SS-101-002-GEOSY	SO	N	12/4/2018	1	22	9		68	
GS-SS-101	GS-SS-101-003-CDM	GS-SS-101-003-GEOSY	SO	N	12/4/2018						51
GS-SS-111	GS-SS-111-001-CDM	GS-SS-111-001-GEOSY	SO	N	12/4/2018	1	22	9		68	
GS-SS-111	GS-SS-111-003-CDM	GS-SS-111-003-GEOSY	SO	N	12/4/2018						51
MW-10	MW-10-CDM	MW-10-GEOSY	WG	N	12/20/2018	1	22	9	21	68	51
MW-17	MW-17-CDM	MW-17-GEOSY	WG	N	12/27/2018	1	22	9	21	68	51
	TB-122718		WQ	ТВ	12/27/2018						51

Notes:

1. Numbers in the table represent the number of compounds reported in each CDM Smith analytical group.

Abbreviations:

ID = identification SVOC = semivolatile organic compound

N = normal TB = trip blank

PCB = polychlorinated biphenyl VOC = volatile organic compound

SO = soil WG = groundwater SOP = standard operating procedure WQ = water quality

ParameterMethod*ParameterMethod*MercuryISM02.4PesticidesSOM02.4

MercuryISM02.4PesticidesSOM02.4MetalsISM02.4SVOCSOM02.4PCBsSOM02.4VOCSOM02.4

*The methods listed are those included in the electronic data deliverables from the laboratories and are based on SOPs consistent with the quality assurance project plan requested methods.



Data Quality Indicators ([OQIs) and Corresponding Quality Control (QC) Parameters					
Data Quality Indicators QC Parameters Evaluated in Data Review/Validation						
Precision	Relative percent difference (RPD) or percent difference (% D) values of:					
	Initial calibration and continuing calibration verifications					
	Laboratory control sample (LCS)/LCS duplicates					
	Matrix spike/matrix spike duplicate (MS/MSD)					
	ICP inter-element interference check samples					
	Serial dilutions (Inductively Coupled Plasma [ICP] metals)					
Accuracy/Bias	Percent Recovery (% R) values of:					
	Initial calibration and continuing calibration verifications					
	LCS/LCSDs					
	MS/MSDs					
	Surrogate spikes					
	Deuterated monitoring compounds (DMC)					
	Laboratory control samples (LCS)					
	Post digestion spikes					
	Results of:					
	Instrument and calibration blanks					
	Method (preparation) blanks					
	Trip blanks					
Representativeness	Results of all blanks					
	Sample integrity (Chain-of Custody and sample receipt forms)					
	Holding times					
	Adequacy of sample dilutions					
	Method blanks					
Comparability	Sample-specific reporting limits (RLs)					
	Sample collection methods					
	Laboratory analytical methods					
	Split samples					
Completeness	Data qualifiers					
	Laboratory deliverables					
	Requested/reported results					
	Field sample collection (primary and QC samples)					
	Contract compliance (i.e., method and instrument QC within limits)					
Sensitivity	Sample RLs					
	Project action limits (PALs)					
	Project quantitation limit goals					



Activity	Comment						
Field Sampling							
Discuss sampling problems and field conditions that affect data usability.	No Field Change Notifications to the Final Quality Assurance Plan (QAPP) (July 2018) were required.						
Are samples representative of receptor exposure for this medium (e.g. sample depth, grab versus composite, filtered versus unfiltered, low flow, etc.)?	Samples were kept at 0 to 6 degrees Celsius and were received intact at the laboratories. In addition, the split sampling design was developed to determine accuracy of the potentially responsible party (PRP)'s investigation results.						
Assess the effect of field QC results on data usability.	CDM Smith did not accept equipment blank split samples during the event – they are not typically required for oversight assignments. Trip Blanks – Acetone was detected in a trip blank just above the contract required quantitation limit (CRQL) (Table 2-3). The trip blank detection required no sample qualifications.						
Summarize the effect of field sampling issues on the data assessment, if applicable.	None.						
	Analytical Techniques						
Were the analytical methods appropriate for quantitative data assessment?	Yes. Analytical methods used for the soil and groundwater samples were appropriate for oversight evaluation. Volatile organic compounds (VOCs), semivolatile organic compounds (SVOCs), pesticides, and polychlorinated biphenyls (PCBs) were analyzed by United States Environmental Protection Agency (EPA) method SOM02.4. EPA analyzed metals and mercury by method ISM02.4. A list of analyses is included on Table 1-1, Summary of Split Samples.						
Were detection limits adequate?	Except for diluted results, sample reporting limits are at the CRQLs which facilitates comparison of the data with the QAPP PALs. Organic and inorganic samples diluted were due to high concentrations. This did not impact the quality of the comparison since generally the PRP results were higher due to dilution as well.						



Activity	Activity Comment					
Analytical Techniques (Continued)						
Summarize the effect of analytical technique issues on	Internal Standards - Internal standards performance criteria ensure that gas chromatography (GC)/ mass spectrometry (MS) sensitivity and response are stable during every analytical run.					
the data assessment, if applicable.	VOCs internal standards results outside criteria are estimated.					
ii applicasie.	Compound Identification - The retention times of reported compounds must fall within the calculated retention time windows for two chromatographic columns and a GC/MS confirmation is required if PCB concentration exceeds 10 nanograms per milliliter (ng/mL) in the final sample extract. Several PCB Aroclors results are estimated.					
	Analytical Blank - Laboratory method blank detections resulted in qualification of VOCs and metals; associated sample results are qualified as non-detect at the CRQL.					
	Other Problems - Several sample results did not meet criteria for other QC problems identified in the data validation (DV) reports - VOC sample results are qualified as estimated. One VOC result is qualified "J" as the reported value is over the calibration range. The dilution analysis was not performed for this VOC sample (GS-B-101-004-CDM).					
	Reconciliation of Data Quality Objectives					
Precision - How were duplicates handled?	Since CDM Smith accepted split samples from the PRP, field duplicate samples were not required.					
	MS/MSD – Several PCB Aroclors and one pesticide MS/MSD do not meet QC criteria. Affected sample results are qualified as estimated.					
	Inductively Coupled Plasma (ICP) Serial Dilution (Inorganics) - Several ICP serial dilutions do not yield acceptable % D. Associated metal results are estimated.					
Accuracy - How were split samples handled?	As part of the oversight assignment, CDM Smith was tasked to accept split samples for 10 percent of the PRP samples by matrix and compare the split sample results with the PRP data. The split samples were subsequently submitted to the Contract Laboratory Program (CLP) laboratory for analyses.					
	CDM Smith accepted soil and groundwater split samples at a rate of 10 percent from the PRP contractor. Split and PRP samples were compared, and results are shown on Table 3s.					
	Splits and PRP soil samples were evaluated using a criterion of 100 percent RPD applied when both results were greater than five times the EPA's CRQL. For all other data sets, the absolute difference (ABS) should be equal to or less than two times the EPA's CRQL.					



Activity	Comment						
	Reconciliation of Data Quality Objectives (Continued)						
Accuracy - How were split samples handled? - Continued	Splits and PRP groundwater samples were evaluated using a criterion of 50 percent RPD applied when both results were greater than five times the EPA's CRQL. For all other data sets the ABS should be equal to or less than two times the EPA's CRQL.						
	Soil - Target Analyte List (TAL) metals, including mercury, Target Compound List (TCL) PCBs, TCL SVOCs, and TCL VOCs.						
	The outliers below failed the applicable RPD or ABS criteria:						
	■ TAL Metals and Hg — RPD/ABS results for 15 analytes exceeded out of 202 calculable pairs						
	 TCL PCBs- RPD/ABS results for six analytes exceeded out of 23 calculable pairs TCL SVOC- RPD/ABS results for 67 analytes exceeded out of 186 calculable pairs TCL VOC- RPD/ABS results for 46 analytes exceeded out of 89 calculable pairs 						
	As shown on Tables 3-1a and 3-1b, most of the soil split results met criteria. Refer to Section 4 of the text for additional data comparison discussion.						
	Groundwater - TAL metals, including mercury, TCL PCBs, TCL pesticides, TCL SVOCs, and TCL VOCs						
	The outliers below failed the applicable RPD or ABS criteria:						
	 TAL Metals and Hg – RPD/ABS results for eight analytes exceeded out of 37 calculable pairs 						
	 TCL Pesticide – ABS result for one analyte exceeded out of one calculable pair TCL VOC – RPD/ABS results for five analytes exceeded out of 29 calculable pairs 						
	As shown on Table 3-2, most of the groundwater split results met criteria. Refer to Section 4 of the text for additional data comparison discussion.						
	DMCs and Surrogate Recoveries – Several DMCs and surrogates exceeded QC criteria. This affected sample results for some VOC, SVOCs, and PCB which are estimated. SVOC 1,4-dioxane was rejected in two samples (GS-B-112-001-CDM and GS-B-114-001-CDM) due to surrogate recoveries.						



Activity	Comment						
1	Reconciliation of Data Quality Objectives (Continued)						
Accuracy - How were split samples handled? - Continued	Percent Relative Standard Deviation (% RSD) and Percent Difference (% D) - Percent RSD is calculated from the initial calibration and used to indicate stability of the specific compound response factor over increasing concentration. % D compares the response factor of the continuing calibration check to the mean relative response factor (RRF) from the initial calibration. % D is a measure of the instrument's daily performance. A value outside of these limits indicates potential detection and quantitation errors. % RSD and % D recoveries outside control limits for some VOC and SVOCs were estimated by the data validator. Matrix Spike – Several MS or MS/MSD sample results do not meet QC criteria. Associated PCBs, pesticides, metals and mercury sample results are qualified as						
Representativeness - Indicate any problems	estimated. As specified in the QAPP, no equipment blanks were collected during this split sampling event.						
associated with data representativeness (e.g., trip blank or rinsate blank contamination, chain of custody problems, etc.).	Trip Blanks – Acetone was detected above the CRQL in the trip blank, but this chemical is not a contaminant of concern (Table 2-3). No validation actions were required; therefore, the data validator did not qualify sample results.						
Completeness - Indicate any problems associated with data completeness (e.g.,	Completeness of the data set achieved by CDM Smith is presented in Table 2-2s which only includes environmental samples, not QC samples. The list of collected samples and analytical parameters are shown on Table 1-1. For data collected and judged to be valid, refer to Tables 2-2a and 2-2b:						
incorrect sample analysis, incomplete sample records,	 Soil results were 99.85 percent complete Groundwater results were 100 percent complete 						
problems with field procedures, etc.).	The overall goal was to generate a complete data set for at least 90 percent of the samples planned for collection and 90 percent valid data from the samples analyzed. Therefore, CDM Smith achieved the data completeness goal in the Final QAPP.						
	The data validation narratives indicate that analyses generally met method QC criteria.						
	One hundred percent of planned TCL/TAL parameter split samples were accepted. The data set is complete based on the 10 percent of planned split samples acceptance goal. During the field event, the PRPs added dissolved metals and mercury analyses. CDM Smith did not accept samples for these parameters since these were not scoped. Per EPA's direction, extractable petroleum hydrocarbons were not collected since the CLP laboratory would not analyze the samples.						



Activity	Comment							
	Data Validation and Interpretation							
Comparability - Indicate any problems associated with data comparability.	In general, the PRP and CDM Smith split data sets are comparable for VOCs, SVOCs, pesticides, PCBs, and metals including mercury. Refer to Section 4 of the text for additional data comparison discussion.							
Were the DQOs specified in the QAPP satisfied?	Yes, the data quality objectives (DQOs) identified in the QAPP were satisfied.							
Summarize the effect of DQO issues on the data assessment, if applicable.	There are no DQO issues that affect data usability.							
What are the data validation requirements?	Page 1 shows the DQIs and Corresponding QC Parameters and the specific topics covered by the data validation review. 100 percent of the data were validated.							
What method or guidance was used to	The data validation criteria in the QAPP were applied: Standard operating procedures (SOPs) HW-33A revision (rev.) 1; HW-34A rev. 1; HW-							
validate the data?	35A rev. 1; HW-36A rev. 1; HW-37A, rev. 0; HW-3A rev. 1; HW-3B rev 1.; HW-3C rev. 1							
Was the data validation method consistent with guidance? Discuss any discrepancies.	Yes. The data validation method was consistent with regional guidance.							
Were all data qualifiers defined? Discuss those which were not.	Yes. All data qualifiers were defined.							
Which qualifiers represent useable data?	J, J+, J-, NJ or JN, U, and UJ							
Which qualifiers represent unusable data?	R							
How are tentatively identified compounds handled?	These data were not evaluated under this study.							



Activity	Comment					
	Data Validation and Interpretation (Continued)					
Summarize the effect of data validation and interpretation issues on the data assessment, if applicable.	Data qualified with an "R" will not be used.					
Additional notes:	None.					

Note: This Worksheet summarizes the data usability analysis and conclusions. Refer to specific pages in the oversight split comparison report for more details on the information presented here.



Table 2-2a CDM Smith Soil Completeness Pierson's Creek Site, Operable Unit 2 Newark, New Jersey

Analytical Parameter	Non- Detects	No. of Hits	No. of Rejects	No. of Estimated Hits	Total	Percent Rejected	Percent Estimated Hits
VOCs	341	31	0	36	408	0.00	8.82
SVOCs	470	63	2	77	612	0.33	12.58
PCBs	67	8	0	6	81	0.00	7.41
Metals	17	158	0	23	198	0.00	11.62
Mercury	0	8	0	1	9	0.00	11.11
Sum	895	268	2	143	1308	0.15	10.93

Percent of all Data Rejected	0.15	
Percent of all SVOC Rejected	0.33	
Percent of all Hits Estimated	10.93	(does not include estimated non-detect data)
Percent Complete (judged valid)	99.85	(Includes all estimated data)

Notes:

The counts and calculations above include only environmental samples; field duplicate, field blank, and trip blanks are not included.

PCB = polychlorinated biphenyl SVOC = semivolatile organic compound VOC = volatile organic compound



Table 2-2b CDM Smith Groundwater Completeness Pierson's Creek Site, Operable Unit 2 Newark, New Jersey

Analytical Parameter	Non- Detects	No. of Hits	No. of Rejects	No. of Estimated Hits	Total	Percent Rejected	Percent Estimated Hits
VOCs	76	18	0	8	102	0.00	7.84
SVOCs	134	1	0	1	136	0.00	0.74
Pesticides	42	0	0	0	42	0.00	0.00
PCBs	18	0	0	0	18	0.00	0.00
Metals	12	27	0	5	44	0.00	11.36
Mercury	1	1	0	0	2	0.00	0.00
Sum	283	47	0	14	344	0.00	4.07

Percent of all Data Rejected	0.00	
Percent of all Hits Estimated	4.07	(does not include estimated non-detect data)
Percent Complete (judged valid)	100	(Includes all estimated data)

Notes:

The counts and calculations above include only environmental samples; field duplicate, field blank, and trip blanks are not included.

PCB = polychlorinated biphenyl SVOC = semivolatile organic compound VOC = volatile organic compound



Table 2-3 Trip Blank Results Pierson's Creek Site, Operable Unit 2 Newark, New Jersey

Chemical Name	CBOL	Unit	TB-122718
Chemical Name	CRQL	Onit	12/27/2018
VOCs			
Acetone	5	μg/L	6

Notes:

- 1. Samples and compounds with no detects are not included in the table above.
- 2. Hits at or above the CRQL are highlighted yellow and bolded.

CRQL = contract required quantitation limits

TB = trip blank

μg/L = microgram per liter



	2 x		GS-B-101-001-CD	M	GS-B-101-001-GEC	OSY			GS-B-102-001-C	M	GS-B-102-001-GEOS	Υ		
Chemical Name	CRQL	Unit	12/14/2018		12/14/2018		RPD	ABS	12/17/2018		12/17/2018		RPD	ABS
	CNQL		GS	-B-1	101-001				GS	S-B-	102-001			
Metals														
Aluminum	40	mg/kg	3320		3120		6.21	NA	8140		7780		4.52	NA
Antimony	2	mg/kg	1.4	J	4.28		NA	2.88	0.93	J	8.88		NA	7.95
Arsenic	1	mg/kg	9.3		16.8		57.5	NA	19.4		46.8		82.8	NA
Barium	10	mg/kg	129		169	J	26.8	NA	154		159		3.19	NA
Beryllium	1	mg/kg	0.32	J	0.267	J	NA	0.05	0.46		0.279 J		NA	0.18
Cadmium	1	mg/kg	1.8		2.13	J	NA	0.33	5.2		7.41 J		35.1	NA
Calcium	1000	mg/kg	6050		6400		5.62	NA	32800		37300		12.8	NA
Chromium	2	mg/kg	19.4		21		7.92	NA	55.8		99.4		56.2	NA
Cobalt	1	mg/kg	5.2		4.88		6.35	NA	40.8		41.5		1.7	NA
Copper	2	mg/kg	126		114		10	NA	195		464		81.6	NA
Iron	20	mg/kg	11300		10400		8.29	NA	24800		36400		37.9	NA
Lead	1	mg/kg	275		289		4.96	NA	462		716		43.1	NA
Magnesium	1000	mg/kg	1010		938		NA	72	3760		5620		39.7	NA
Manganese	1	mg/kg	80.9		105		25.9	NA	197		455		79.1	NA
Mercury	0	mg/kg	30.7		20.8		38.4	NA	641		445		36.1	NA
Nickel	1	mg/kg	19.4		20.4		5.03	NA	56.9		110 J-	-	63.6	NA
Potassium	1000	mg/kg	415	J	255		NA	160	758		768		NA	10
Selenium	5	mg/kg	2.9	J	1.66	U	NC	NC	3	UJ	2.26 J		NA	0.74
Silver	1	mg/kg	0.83	J	1.48		NA	0.65	7.3	J-	11.8		47.1	NA
Sodium	1000	mg/kg	629		673		NA	44	1030		947 J-	+	NA	83
Thallium	1	mg/kg	0.67	J	1.66	U	NA	0.99	1.3	J	1.86 U	J	NA	0.56
Vanadium	5	mg/kg	23.3		21.2		9.44	NA	35.9		49.9 J-	-	32.6	NA
Zinc	2	mg/kg	205		222		7.96	NA	3530		7210		68.5	NA
Polychlorinated Biphenyl														
Aroclor 1016	66	μg/kg	37	U	30	С	NC	NC	40	U	190 U	J	NC	NC
Aroclor 1221	66	μg/kg	37	U	30	U	NC	NC	40	U	190 U	J	NC	NC
Aroclor 1232	66	μg/kg	37	U	30	U	NC	NC	40	U	190 U	J	NC	NC
Aroclor 1242	66	μg/kg	37	U	352	J	NA	315	40	U	190 U	J	NC	NC
Aroclor 1248	66	μg/kg	37	U	30	U	NC	NC	40	U	190 U	J	NC	NC
Aroclor 1254	66	μg/kg	460		532	J	14.5	NA	580		1570 J		92.1	NA
Aroclor 1260	66	μg/kg	220	NJ	207	J	6.09	NA	330	J	418 J		23.5	NA
Aroclor 1262	66	μg/kg	37	U	30	U	NC	NC	40	U	190 U	J	NC	NC



	2 x		GS-B-101-001-CD	M	GS-B-101-001-GEC	DSY			GS-B-102-001-CD	М	GS-B-102-001-GEOS	SY		
Chemical Name	CRQL	Unit	12/14/2018		12/14/2018		RPD	ABS	12/17/2018		12/17/2018		RPD	ABS
	CKQL		GS	-B-1	L01-001				GS-	-B-10	02-001			
Aroclor 1268	66	μg/kg	37	U	30	С	NC	NC	40 l	U	190 Լ	U	NC	NC
Semivolatile Organic Compound														
1,1'-Biphenyl	340	μg/kg	190	U	410	С	NC	NC	200 Լ	U	440 l	U	NC	NC
1,2,4,5-Tetrachlorobenzene	340	μg/kg	190	U	180	U	NC	NC	200 l	J	200 Լ	U	NC	NC
1,4-Dioxane ²	340	μg/kg	75	U			NC	NC	81 l	J			NC	NC
2,2'-Oxybis(1-chloropropane)	340	μg/kg	370	U	210	UJ	NC	NC	400 l	U	230 Լ	UJ	NC	NC
2,3,4,6-Tetrachlorophenol	134	μg/kg	190	U	180	UJ	NC	NC	200 l	J	200 F	R	NC	NC
2,4,5-Trichlorophenol	340	μg/kg	190	U	180	UJ	NC	NC	200 Լ	U	200 Լ	UJ	NC	NC
2,4,6-Trichlorophenol	340	μg/kg	190	U	100	UJ	NC	NC	200 l	U	100 Լ	UJ	NC	NC
2,4-Dichlorophenol	340	μg/kg	190	U	80	U	NC	NC	200 l	J	90 ا	UJ	NC	NC
2,4-Dimethylphenol	340	μg/kg	190	U	170	J	NC	NC	200 l	U	180 l	UJ	NC	NC
2,4-Dinitrophenol	340	μg/kg	370	UJ	250	R	NC	NC	400 l	IJ	280 F	R	NC	NC
2,4-Dinitrotoluene	660	μg/kg	190	U	90	J	NC	NC	200 l	J	90 l	UJ	NC	NC
2,6-Dinitrotoluene	340	μg/kg	190	U	70	U	NC	NC	200 l	J	70 l	UJ	NC	NC
2-Chloronaphthalene	340	μg/kg	190	U	180	U	NC	NC	200 Լ	J	200 l	UJ	NC	NC
2-Chlorophenol	340	μg/kg	190	U	50	UJ	NC	NC	200 Լ	J	60 l	UJ	NC	NC
2-Methylnaphthalene	340	μg/kg	190	U	92	J	NA	98	200 Լ	J	52 J	J	NA	148
2-Methylphenol	340	μg/kg	370	U	180	U	NC	NC	400 l	J	200 l	UJ	NC	NC
2-Nitroaniline	660	μg/kg	190	UJ	180	U	NC	NC	200 Լ	IJ	200 Լ	U	NC	NC
2-Nitrophenol	340	μg/kg	190	U	380	IJ	NC	NC	200 l	J	410 F	R	NC	NC
3,3'-Dichlorobenzidine	340	μg/kg	370	U	140	UJ	NC	NC	400 l	J	150 Լ	UJ	NC	NC
3-Nitroaniline	660	μg/kg	370	UJ	180	UJ	NC	NC	400 l	IJ	200 Լ	UJ	NC	NC
4,6-Dinitro-2-methylphenol	660	μg/kg	370	UJ	250	R	NC	NC	400 l	IJ	280 F	R	NC	NC
4-Bromophenyl-phenylether	660	μg/kg	190	U	180	U	NC	NC	200 l	J	200 l	U	NC	NC
4-Chloro-3-methylphenol	340	μg/kg	190	U	180	U	NC	NC	200 Լ	J	200 Լ	U	NC	NC
4-Chloroaniline	340	μg/kg	370		180	U	NC	NC	400 l	J	200 l	U	NC	NC
4-Chlorophenyl-phenylether	660	μg/kg	190	U	180	U	NC	NC	200 Լ	J	200 Լ	U	NC	NC
4-Methylphenol	340	μg/kg	370				NC	NC	400 l	J			NC	NC
4-Nitroaniline	660	μg/kg	370	UJ	180	U	NC	NC	400 l	IJ	200 Լ	_	NC	NC
4-Nitrophenol	660	μg/kg	370	UJ	250	UJ	NC	NC	400 l	IJ	270 F	R	NC	NC
Acenaphthene	660	μg/kg	190		53		NA	137	87 J		150 J		NA	63
Acenaphthylene	340	μg/kg	190	U	28	J	NA	162	200 l	IJ	50 J	J	NA	150



	2 x		GS-B-101-001-CD	M	GS-B-101-001-GEOSY	1		GS-B-102-001-CDM	GS-B-102-001-GEOSY	,	
Chemical Name	CRQL	Unit	12/14/2018		12/14/2018	RP	O AB	12/17/2018	12/17/2018	RPD	ABS
	CNQL		GS-	-B-1	101-001			GS-B-	-102-001		
Acetophenone	340	μg/kg	370		180 U	N	C NO		200 U	NC	NC
Anthracene	660	μg/kg	190		78 J	N.			440	NA	190
Atrazine	340	μg/kg	370	U	150 U	N	C NO	400 U	150 U.	J NC	NC
Benzaldehyde	660	μg/kg	370	U	230 U	N	C NO	400 U	250 U	NC	NC
Benzo(a)anthracene	660	μg/kg	91	J	160 J	N.	4 69	890	1400 J	NA	510
Benzo(a)pyrene	340	μg/kg	93	J	160 J	N.	67	880	1700 J	63.6	NA
Benzo(b)fluoranthene	340	μg/kg	130	J	250 J	N.	120	1100	1900 J	53.3	NA
Benzo(g,h,i)perylene	340	μg/kg	50	J	120 J	N.	70	530	1200 J	NA	670
Benzo(k)fluoranthene	340	μg/kg	49	J	79 J	N.	30	350	590 J	NA	240
Bis(2-chloroethoxy)methane	340	μg/kg	190	U	200 UJ	N	C NO	200 U	200 U	NC	NC
Bis(2-chloroethyl)ether	340	μg/kg	370	U	60 UJ	N	C NO	400 U	70 UJ	I NC	NC
Bis(2-ethylhexyl)phthalate	340	μg/kg	250		2600 J	N.	A 235	<mark>0</mark> 280	640	NA	360
Butylbenzylphthalate	660	μg/kg	190	U	180 U	N	C NO	200 U	200 U	NC	NC
Caprolactam	340	μg/kg	370	U	180 U	N	C NO	400 U	200 U	NC	NC
Carbazole	660	μg/kg	370	כ	33 J	N.	337	7 59 J	140 J	NA	81
Chrysene	660	μg/kg	99	J	210	N.	11:	920	1400 J	NA	480
Dibenzo(a,h)anthracene	660	μg/kg	190	J	34 J	N.	150	160 J	260	NA	100
Dibenzofuran	340	μg/kg	190	J	36 J	N.	154	200 U	66 J	NA	134
Diethylphthalate	340	μg/kg	190	U	180 U	N	C NO	200 U	200 U	NC	NC
Dimethylphthalate	340	μg/kg	100	J	180 U	N.	A 80	140 J	200 U	NA	60
Di-n-butylphthalate	340	μg/kg	190	J	160 J	N.	30	170 J	56 J	NA	114
Di-n-octylphthalate	340	μg/kg	370	U	180 U	N	C NO	400 U	200 U	NC	NC
Fluoranthene	660	μg/kg	190	J	420	N.	230	1300	3000 J	NA	1700
Fluorene	660	μg/kg	190	U	78 J	N.	112	2 87 J	170 J	NA	83
Hexachlorobenzene	340	μg/kg	190	U	50 UJ	N	C NO	200 U	50 U	NC	NC
Hexachlorobutadiene	340	μg/kg	190	J	60 U	N	C NO	200 U	70 UJ	I NC	NC
Hexachlorocyclopentadiene	340	μg/kg	370	U	510 R	N	C NO	400 U	560 R	NC	NC
Hexachloroethane	660	μg/kg	190	U	80 U	N	C NO	200 U	90 U	NC	NC
Indeno(1,2,3-cd)pyrene	340	μg/kg	48	J	130 J	N.	82	490	1200 J	NA	710
Isophorone	340	μg/kg	190	U	60 UJ	N	C NC	200 U	60 U	NC	NC
Naphthalene	340	μg/kg	410		1500 J	N.	109	0 200 U	100 J	NA	100
Nitrobenzene	340	μg/kg	190	U	80 U	N	C NO	200 U	80 U	NC	NC
N-Nitroso-di-n-propylamine	340	μg/kg	190	U	40 U	N	C NO	200 U	40 U	NC	NC



Table 3-1a

Soil Samples Data Comparison - Metals, PCBs, and SVOCs Pierson's Creek Site, Operable Unit 2 Newark, New Jersey

2)			GS-B-101-001-CDM	GS-B-101-001-GEOSY			GS-B-102-001-CDN	GS-B-102-001-GEOSY		
Chemical Name	CRQL	Unit	12/14/2018	12/14/2018	RPD	ABS	12/17/2018	12/17/2018	RPD	ABS
	CNQL		GS-B-	101-001			GS-I	3-102-001		
N-Nitrosodiphenylamine	340	μg/kg	230	1000 J	NA	770	200 U	40 U	NC	NC
Pentachlorophenol	660	μg/kg	370 U	110 UJ	NC	NC	400 U	120 UJ	NC	NC
Phenanthrene	340	μg/kg	130 J	330 J	NA	200	1000	1500 J	40	NA
Phenol	660	μg/kg	99 J	180 U	NA	81	130 J	200 UJ	NA	70
Pyrene	340	μg/kg	200	360	NA	160	1900	3300 J	53.8	NA

Notes:

- 1. Sample pairs with RPD or ABS outside of criteria in the table are highlighted in yellow
- 2. PRP's laboratory analyzed 1,4-dioxane as a volatile organic compound.

ABS = absolute difference

CRQL = contract required quantitation limit

mg/kg = milligram per kilogram

NA = not applicable

NC = not calculable

PCB = polychlorinated biphenyl

RPD = relative percent difference

SVOC = semivolatile organic compound

μg/kg = microgram per kilogram

Data Validation Qualifiers

J = The analyte was positively identified, and the associated numerical value is the approximate concentration of the analyte in the sample.

J+ = The result is an estimated quantity, but the result may be biased high.

J- = The result is an estimated quantity, but the result may be biased low.

NJ = The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents its approximate concentration.

R = The sample results are unusable due to poor data quality; certain criteria were not met. The analyte may or may not be present in the sample at the reported concentration.

U = The analyte was not detected at a level greater than or equal to the level of the sample CRQL.

UJ = The analyte was not detected; however, the reported CRQL is approximate and may be inaccurate or imprecise.



	2 x		GS-B-102-002-CDN	Λ	GS-B-102-002-GEC	DSY				M	GS-B-111-001-GEOS	SY		
Chemical Name	CRQL	Unit	12/17/2018		12/17/2018		RPD	ABS	12/17/2018		12/17/2018		RPD	ABS
	ChQL		GS-E	3-10	02-002				GS	S-B-	111-001			
Metals														
Aluminum	40	mg/kg	17000		10400		48.2	NA	8500		8930		4.93	NA
Antimony	2	mg/kg	0.53	J	0.859	J	NA	0.33	4.7	J	4.53 l	U	NC	NC
Arsenic	1	mg/kg	8		4.32		59.7	NA	2.3		3.32		NA	1.02
Barium	10	mg/kg	364		161	J	77.3	NA	63.3		44.4		35.1	NA
Beryllium	1	mg/kg	0.74		0.465	J	NA	0.28	1.3		0.672		NA	0.63
Cadmium	1	mg/kg	0.4	J	1.19	U	NA	0.79	0.75		0.273 J	J	NA	0.48
Calcium	1000	mg/kg	6340		3400	J	60.4	NA	14900		20200		30.2	NA
Chromium	2	mg/kg	15.9		9.14	J	54	NA	23.2		25.5		9.45	NA
Cobalt	1	mg/kg	8.6		3.26		90.1	NA	34.8		27.6		23.1	NA
Copper	2	mg/kg	34.3		20.3		51.3	NA	175		120		37.3	NA
Iron	20	mg/kg	11500		6020		62.6	NA	14700		15300		4	NA
Lead	1	mg/kg	60.9		21.9		94.2	NA	248		175		34.5	NA
Magnesium	1000	mg/kg	493	J	225	J	NA	268	5430		6170		12.8	NA
Manganese	1	mg/kg	50.9		24.7		69.3	NA	106		118		10.7	NA
Mercury	0	mg/kg	12.5		9.77		24.5	NA	74.5		75.6		1.47	NA
Nickel	1	mg/kg	12.1		7.83	J	42.8	NA	80.6		48.7		49.3	NA
Potassium	1000	mg/kg	1210		796		NA	414	466		775		NA	309
Selenium	5	mg/kg	4.2	U	0.537	J	NA	3.66	2.7	J	1.82 l	U	NC	NC
Silver	1	mg/kg	1.2	U	1.19	J	NC	NC	0.83		2.21		NA	1.38
Sodium	1000	mg/kg	2750		1680	J	NA	1070	922		1010		NA	88
Thallium	1	mg/kg	0.56	J	2.39	U	NA	1.83	1.1	J	1.82 l	U	NA	0.72
Vanadium	5	mg/kg	21.3		12.2		NA	9.1	31.7		41.2		26.1	NA
Zinc	2	mg/kg	74.5		26.7		94.5	NA	384		359		6.73	NA
Polychlorinated Biphenyl														
Aroclor 1016	66	μg/kg	56	U	50	U	NC	NC	37	U	30 l	U	NC	NC
Aroclor 1221	66	μg/kg	56	U	50	U	NC	NC	37	U	30 L	U	NC	NC
Aroclor 1232	66	μg/kg	56	U	50	U	NC	NC	37	U	30 l	U	NC	NC
Aroclor 1242	66	μg/kg	56	U	50	U	NC	NC	37	U	30 l	U	NC	NC
Aroclor 1248	66	μg/kg	56	U	48.4	J	NA	7.6	37	U	30 l	U	NC	NC
Aroclor 1254	66	μg/kg	56	U	66.7	J	NA	10.7	37	U	52.3 J	J	NA	15.3
Aroclor 1260	66	μg/kg	56	U	29.9	J	NA	26.1	37	U	32.2 J	J	NA	4.8
Aroclor 1262	66	μg/kg	56	U	50	U	NC	NC	37	U	30 l	U	NC	NC



	2 x		GS-B-102-002-CDN	Л	GS-B-102-002-GEO	SY			GS-B-111-001-CD	M	GS-B-111-001-GEOSY	,		
Chemical Name	CRQL	Unit	12/17/2018		12/17/2018		RPD	ABS	12/17/2018		12/17/2018	RP) A	ABS
	CITQL			_	02-002						111-001		\perp	
Aroclor 1268	66	μg/kg	56	U	50	U	NC	NC	37	U	30 U	NO		NC
Semivolatile Organic Compound														
1,1'-Biphenyl	340	μg/kg	290	U	560	U	NC	NC	190	U	460 U	N	1	NC
1,2,4,5-Tetrachlorobenzene	340	μg/kg	290	U	250	U	NC	NC	190	U	200 U	N	. 1	NC
1,4-Dioxane ²	340	μg/kg	110	U			NC	NC	75	U		NO	1	NC
2,2'-Oxybis(1-chloropropane)	340	μg/kg	560	U	300	U	NC	NC	370	U	230 U	N	: 1	NC
2,3,4,6-Tetrachlorophenol	134	μg/kg	290	U	250	U	NC	NC	190	U	200 UJ	NO	. 1	NC
2,4,5-Trichlorophenol	340	μg/kg	290	U	250	U	NC	NC	190	U	200 UJ	NO	. 1	NC
2,4,6-Trichlorophenol	340	μg/kg	290	U	140	U	NC	NC	190	U	100 UJ	NO	. 1	NC
2,4-Dichlorophenol	340	μg/kg	290	U	110	U	NC	NC	190	U	90 UJ	NO	1	NC
2,4-Dimethylphenol	340	μg/kg	290	U	230	U	NC	NC	190	U	180 UJ	NO	. 1	NC
2,4-Dinitrophenol	340	μg/kg	560	U	360	U	NC	NC	370	UJ	280 UJ	N	. 1	NC
2,4-Dinitrotoluene	660	μg/kg	290	U	110	U	NC	NC	190	U	100 U	NO	1	NC
2,6-Dinitrotoluene	340	μg/kg	290	U	90	U	NC	NC	190	U	70 U	NO	. 1	NC
2-Chloronaphthalene	340	μg/kg	290	U	250	U	NC	NC	190	U	200 U	NO	1	NC
2-Chlorophenol	340	μg/kg	290	U	80	U	NC	NC	190	U	60 UJ	NO	1	NC
2-Methylnaphthalene	340	μg/kg	290	U	300	U	NC	NC	190	U	230 U	NO	1	NC
2-Methylphenol	340	μg/kg	560	U	250	U	NC	NC	370	U	200 UJ	NO	1	NC
2-Nitroaniline	660	μg/kg	290	U	250	U	NC	NC	190	UJ	200 U	NO	1	NC
2-Nitrophenol	340	μg/kg	290	U	540	U	NC	NC	190	U	430 UJ	NO	1	NC
3,3'-Dichlorobenzidine	340	μg/kg	560	U	180	UJ	NC	NC	370	U	150 UJ	NO	1	NC
3-Nitroaniline	660	μg/kg	560	U	250	UJ	NC	NC	370	UJ	200 UJ	NO	1	NC
4,6-Dinitro-2-methylphenol	660	μg/kg	560	U	360	U	NC	NC	370	UJ	280 U	NO	1	NC
4-Bromophenyl-phenylether	660	μg/kg	290	U	250	U	NC	NC	190	U	200 U	NO	1	NC
4-Chloro-3-methylphenol	340	μg/kg	290	U	250	U	NC	NC	190	U	200 U	NO	. 1	NC
4-Chloroaniline	340	μg/kg	560	U	250	U	NC	NC	370	U	200 U	NO	1	NC
4-Chlorophenyl-phenylether	660	μg/kg	290	U	250	U	NC	NC	190	U	200 U	NO	1	NC
4-Methylphenol	340	μg/kg	560	U			NC	NC	370	U		N	1	NC
4-Nitroaniline	660	μg/kg	560	U	250	U	NC	NC	370	UJ	200 U	N	. 1	NC
4-Nitrophenol	660	μg/kg	560	U	340	U	NC	NC	370	UJ	280 UJ	NO	1	NC
Acenaphthene	660	μg/kg	290		83 .	J	NA	207	190		150 U	N		NC
Acenaphthylene	340	μg/kg	290	U	200	U	NC	NC	190	U	150 U	NO	1	NC



	2 x		GS-B-102-002-CDM	GS-B-102-002-GEOS	Υ		GS-B-111-001-CDM	GS-B-111-001-GEOSY		
Chemical Name	CRQL	Unit	12/17/2018	12/17/2018	RPD	ABS	12/17/2018	12/17/2018	RPD	ABS
	CNQL		GS-B-1	02-002			GS-B-	-111-001		
Acetophenone	340	μg/kg	560 U	250 U	NC	NC	370 U	200 U	NC	NC
Anthracene	660	μg/kg	290 U	150 J	NA	140	190 U	23 J	NA	167
Atrazine	340	μg/kg	560 U	200 U	NC	NC	370 U	150 U	NC	NC
Benzaldehyde	660	μg/kg	560 U	330 U	NC	NC	370 U	250 U	NC	NC
Benzo(a)anthracene	660	μg/kg	130 J	790	NA	660	41 J	70	NA	29
Benzo(a)pyrene	340	μg/kg	210 J	1400	NA	1190	46 J	83 J	NA	37
Benzo(b)fluoranthene	340	μg/kg	260 J	1800	NA	1540	54 J	100	NA	46
Benzo(g,h,i)perylene	340	μg/kg	140 J	1300	NA	1160	190 U	66 J	NA	124
Benzo(k)fluoranthene	340	μg/kg	100 J	530	NA	430	190 U	38 J	NA	152
Bis(2-chloroethoxy)methane	340	μg/kg	290 U	270 U	NC	NC	190 U	210 U	NC	NC
Bis(2-chloroethyl)ether	340	μg/kg	560 U	90 U	NC	NC	370 U	70 U	NC	NC
Bis(2-ethylhexyl)phthalate	340	μg/kg	290 U	250 U	NC	NC	850	290	NA	560
Butylbenzylphthalate	660	μg/kg	290 U	250 U	NC	NC	190 U	200 U	NC	NC
Caprolactam	340	μg/kg	560 U	250 U	NC	NC	370 U	200 U	NC	NC
Carbazole	660	μg/kg	560 U	80 J	NA	480	370 U	200 U	NC	NC
Chrysene	660	μg/kg	130 J	820	NA	690	38 J	74 J	NA	36
Dibenzo(a,h)anthracene	660	μg/kg	290 U	230	NA	60	190 U	70 U	NC	NC
Dibenzofuran	340	μg/kg	290 U	34 J	NA	256	190 U	200 U	NC	NC
Diethylphthalate	340	μg/kg	290 U	250 U	NC	NC	190 U	200 U	NC	NC
Dimethylphthalate	340	μg/kg	130 J	250 U	NA	120	99 J	200 U	NA	101
Di-n-butylphthalate	340	μg/kg	290 U	250 U	NC	NC	190 U	200 U	NC	NC
Di-n-octylphthalate	340	μg/kg	560 U	250 U	NC	NC	370 U	200 U	NC	NC
Fluoranthene	660	μg/kg	170 J	1000 J	NA	830	370 U	100 J	NA	270
Fluorene	660	μg/kg	290 U	250 U	NC	NC	190 U	200 U	NC	NC
Hexachlorobenzene	340	μg/kg	290 U	70 U	NC	NC	190 U	50 U	NC	NC
Hexachlorobutadiene	340	μg/kg	290 U	90 U	NC	NC	190 U	70 U	NC	NC
Hexachlorocyclopentadiene	340	μg/kg	560 U	700 U	NC	NC	370 U	560 U	NC	NC
Hexachloroethane	660	μg/kg	290 U	110 U	NC	NC	190 U	90 U	NC	NC
Indeno(1,2,3-cd)pyrene	340	μg/kg	130 J	1200	NA	1070	190 U	63 J	NA	127
Isophorone	340	μg/kg	290 U	80 U	NC	NC	190 U	60 U	NC	NC
Naphthalene	340	μg/kg	290 U	58 J	NA	232	190 U	200 U	NC	NC
Nitrobenzene	340	μg/kg	290 U	100 U	NC	NC	190 U	80 U	NC	NC
N-Nitroso-di-n-propylamine	340	μg/kg	290 U	60 U	NC	NC	190 U	50 U	NC	NC



Table 3-1a

Soil Samples Data Comparison - Metals, PCBs, and SVOCs Pierson's Creek Site, Operable Unit 2 Newark, New Jersey

	2 x		GS-B-102-002-CDM	GS-B-102-002-GEOSY			GS-B-111-001-CDM	GS-B-111-001-GEOSY		
Chemical Name	CRQL	Unit	12/17/2018	12/17/2018	RPD	ABS	12/17/2018	12/17/2018	RPD	ABS
	CKQL		GS-B-1	.02-002			GS-B-	111-001		
N-Nitrosodiphenylamine	340	μg/kg	290 U	60 U	NC	NC	190 U	40 U	NC	NC
Pentachlorophenol	660	μg/kg	560 U	150 U	NC	NC	370 U	120 UJ	NC	NC
Phenanthrene	340	μg/kg	81 J	560 J	NA	479	190 U	69 J	NA	121
Phenol	660	μg/kg	170 J	250 U	NA	80	98 J	200 UJ	NA	102
Pyrene	340	μg/kg	170 J	990 J	NA	820	190 U	100 J	NA	90

Notes:

- 1. Sample pairs with RPD or ABS outside of criteria in the table are highlighted in yellow
- 2. PRP's laboratory analyzed 1,4-dioxane as a volatile organic compound.

ABS = absolute difference

CRQL = contract required quantitation limit

mg/kg = milligram per kilogram

NA = not applicable

NC = not calculable

PCB = polychlorinated biphenyl

RPD = relative percent difference

SVOC = semivolatile organic compound

μg/kg = microgram per kilogram

Data Validation Qualifiers

J = The analyte was positively identified, and the associated numerical value is the approximate concentration of the analyte in the sample.

J+ = The result is an estimated quantity, but the result may be biased high.

J- = The result is an estimated quantity, but the result may be biased low.

NJ = The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents its approximate concentration.

R = The sample results are unusable due to poor data quality; certain criteria were not met. The analyte may or may not be present in the sample at the reported concentration.

U = The analyte was not detected at a level greater than or equal to the level of the sample CRQL.

UJ = The analyte was not detected; however, the reported CRQL is approximate and may be inaccurate or imprecise.



	2 x		GS-B-112-001-CDN	Λ	GS-B-112-001-GE0	OSY			GS-SS-101-001-CD	M	GS-SS-101-001-GEOSY		
Chemical Name	CRQL	Unit	12/13/2018		12/13/2018		RPD	ABS	12/4/2018		12/4/2018	RPD	ABS
	CKQL		GS-B	3-1:	12-001				GS-S	SS-1	101-001		
Metals													
Aluminum	40	mg/kg	6110		5550		9.61	NA	6270		6460	2.99	NA
Antimony	2	mg/kg	0.44 J		2.02	J	NA	1.58	1.6	J	4.3 J	NA	2.7
Arsenic	1	mg/kg	6.2		6.75	J	8.49	NA	11.5		13.7	17.5	NA
Barium	10	mg/kg	67.4		56.1	J	18.3	NA	142		152	6.8	NA
Beryllium	1	mg/kg	0.32 J		0.193	J	NA	0.13	0.52		0.339 J	NA	0.18
Cadmium	1	mg/kg	3.8		4.18		9.52	NA	1.9		1.58	NA	0.32
Calcium	1000	mg/kg	7680		6720	J	13.3	NA	12300		9120	29.7	NA
Chromium	2	mg/kg	14.3		14		2.12	NA	29		33.4 J	14.1	NA
Cobalt	1	mg/kg	17.8		17.8		0	NA	24.2		30.4	22.7	NA
Copper	2	mg/kg	109		94.8		13.9	NA	121		111	8.62	NA
Iron	20	mg/kg	14900		14500		2.72	NA	25400		20300	22.3	NA
Lead	1	mg/kg	150		131		13.5	NA	805		670 J	18.3	NA
Magnesium	1000	mg/kg	3200		3390		5.77	NA	2050		2760 J	NA	710
Manganese	1	mg/kg	188		205		8.65	NA	300		337	11.6	NA
Mercury	0	mg/kg	61.7		75.4		20	NA	138		157	12.9	NA
Nickel	1	mg/kg	19.5		18.2		6.9	NA	32.5		29.2 J	10.7	NA
Potassium	1000	mg/kg	460		440		NA	20	653		696	NA	43
Selenium	5	mg/kg	2.8 l	J	1.75	J	NC	NC	3 (U	0.578 J	NA	2.42
Silver	1	mg/kg	3		3.1	J	3.28	NA	1.8		2.45	NA	0.65
Sodium	1000	mg/kg	439		382		NA	57	833		875 J+	NA	42
Thallium	1	mg/kg	0.88 J		1.75	J	NA	0.87	1.2	J	1.83 U	NA	0.63
Vanadium	5	mg/kg	34.6		34.7		0.29	NA	26.8		27.4	2.21	NA
Zinc	2	mg/kg	149		138	J	7.67	NA	305		277	9.62	NA
Polychlorinated Biphenyl													
Aroclor 1016	66	μg/kg	36 L	J	30	U	NC	NC	40	U	30 U	NC	NC
Aroclor 1221	66	μg/kg	36 L	J	30	U	NC	NC	40	U	30 U	NC	NC
Aroclor 1232	66	μg/kg	36 L	J	30	U	NC	NC	40	U	30 U	NC	NC
Aroclor 1242	66	μg/kg	36 L	J	30	U	NC	NC	40	U	82.8 JN	NA	42.8
Aroclor 1248	66	μg/kg	36 L	J	30	U	NC	NC	40	U	30 U	NC	NC
Aroclor 1254	66	μg/kg	170		198	J	15.2	NA	180		601 J	108	NA
Aroclor 1260	66	μg/kg	95 N	ΛJ	50.4	J	NA	44.6	170		389 J	78.4	NA
Aroclor 1262	66	μg/kg	36 L	J	30	U	NC	NC	40	U	30 U	NC	NC



	2 x		GS-B-112-001-CD	M	GS-B-112-001-GE	OSY			GS-SS-101-001-CDM	GS-SS-101-001-GEOS	Y		
Chemical Name	CRQL	Unit	12/13/2018		12/13/2018		RPD	ABS	12/4/2018	12/4/2018		RPD	ABS
	CKQL		GS-	-B-1	12-001				GS-SS-	101-001			
Aroclor 1268	66	μg/kg	36	U	30	U	NC	NC	40 U	30 U	J	NC	NC
Semivolatile Organic Compound													
1,1'-Biphenyl	340	μg/kg	930	U	2000	U	NC	NC	210 U	890 U	IJ	NC	NC
1,2,4,5-Tetrachlorobenzene	340	μg/kg	930	U	890	U	NC	NC	210 U	380 U	IJ	NC	NC
1,4-Dioxane ²	340	μg/kg	370	R			NC	NC	81 UJ			NC	NC
2,2'-Oxybis(1-chloropropane)	340	μg/kg	1800	U	1100	U	NC	NC	400 U	460 U	IJ	NC	NC
2,3,4,6-Tetrachlorophenol	134	μg/kg	930	U	890	UJ	NC	NC	210 U	380 U	IJ	NC	NC
2,4,5-Trichlorophenol	340	μg/kg	930	U	890	U	NC	NC	210 U	380 U	IJ	NC	NC
2,4,6-Trichlorophenol	340	μg/kg	930	U	500	U	NC	NC	210 U	210 U	IJ	NC	NC
2,4-Dichlorophenol	340	μg/kg	930	U	430	U	NC	NC	210 U	180 U	IJ	NC	NC
2,4-Dimethylphenol	340	μg/kg	930	U	860	U	NC	NC	210 U	370 U	IJ	NC	NC
2,4-Dinitrophenol	340	μg/kg	1800	UJ	1290	U	NC	NC	400 UJ	560 U	J	NC	NC
2,4-Dinitrotoluene	660	μg/kg	930	U	440	U	NC	NC	210 U	200 U	IJ	NC	NC
2,6-Dinitrotoluene	340	μg/kg	930	U	360	U	NC	NC	210 U	150 U	IJ	NC	NC
2-Chloronaphthalene	340	μg/kg	930	U	890	U	NC	NC	210 U	380 U	IJ	NC	NC
2-Chlorophenol	340	μg/kg	930	U	300	U	NC	NC	210 U	120 U	IJ	NC	NC
2-Methylnaphthalene	340	μg/kg	930	U	1100	U	NC	NC	210 U	130 J		NA	80
2-Methylphenol	340	μg/kg	1800	U	890	U	NC	NC	400 U	380 U	IJ	NC	NC
2-Nitroaniline	660	μg/kg	930	UJ	890	U	NC	NC	210 U	380 U	IJ	NC	NC
2-Nitrophenol	340	μg/kg	930	U	1890	U	NC	NC	210 U	850 U	IJ	NC	NC
3,3'-Dichlorobenzidine	340	μg/kg	1800	U	680	UJ	NC	NC	400 U	300 U	IJ	NC	NC
3-Nitroaniline	660	μg/kg	1800	UJ	890	UJ	NC	NC	400 U	380 U	IJ	NC	NC
4,6-Dinitro-2-methylphenol	660	μg/kg	1800	UJ	1290	U	NC	NC	400 UJ	560 U	IJ	NC	NC
4-Bromophenyl-phenylether	660	μg/kg	930	U	890	UJ	NC	NC	210 U	380 U	IJ	NC	NC
4-Chloro-3-methylphenol	340	μg/kg	930	U	890	U	NC	NC	210 U	380 U	IJ	NC	NC
4-Chloroaniline	340	μg/kg	1800	U	890	U	NC	NC	400 U	380 U	IJ	NC	NC
4-Chlorophenyl-phenylether	660	μg/kg	930	U	890	UJ	NC	NC	210 U	380 U	IJ	NC	NC
4-Methylphenol	340	μg/kg	1800				NC	NC	400 U			NC	NC
4-Nitroaniline	660	μg/kg	1800	_	890	U	NC	NC	400 U	380 U	IJ	NC	NC
4-Nitrophenol	660	μg/kg	1800	UJ	1290	U	NC	NC	400 U	550 U	J	NC	NC
Acenaphthene	660	μg/kg	190		120	-	NA	70	130 J	450 J		NA	320
Acenaphthylene	340	μg/kg	930	U	720	U	NC	NC	84 J	400 J		NA	316



	2 x		GS-B-112-001-CDM	GS-B-112-001-GE0	DSY			GS-SS-101-001-CDM	GS-SS-101-001-GEOSY		
Chemical Name	CRQL	Unit	12/13/2018	12/13/2018		RPD	ABS	12/4/2018	12/4/2018	RPD	ABS
	CNQL		GS-B-1	.12-001				GS-SS-	101-001		
Acetophenone	340	μg/kg	1800 U	890	U	NC	NC	400 U	380 UJ	NC	NC
Anthracene	660	μg/kg	820 J	760		NA	60	550	1100 J	NA	550
Atrazine	340	μg/kg	1800 U	720	U	NC	NC	400 U	310 UJ	NC	NC
Benzaldehyde	660	μg/kg	1800 U	1200	U	NC	NC	400 U	100 J	NA	300
Benzo(a)anthracene	660	μg/kg	2900	2100		32	NA	1200	3900 J	NA	2700
Benzo(a)pyrene	340	μg/kg	2200	1800		20	NA	1000	4100 J	122	NA
Benzo(b)fluoranthene	340	μg/kg	3400	2600		26.7	NA	1400	5200 J	115	NA
Benzo(g,h,i)perylene	340	μg/kg	1000	1000		0	NA	590	2700 J	NA	2110
Benzo(k)fluoranthene	340	μg/kg	1400	1200		15.4	NA	430	1700 J	NA	1270
Bis(2-chloroethoxy)methane	340	μg/kg	930 U	970	IJ	NC	NC	210 U	410 UJ	NC	NC
Bis(2-chloroethyl)ether	340	μg/kg	1800 U	310	IJ	NC	NC	400 U	140 UJ	NC	NC
Bis(2-ethylhexyl)phthalate	340	μg/kg	930 U	120	J	NA	810	210 U	380 UJ	NC	NC
Butylbenzylphthalate	660	μg/kg	930 U	890	כ	NC	NC	210 U	380 UJ	NC	NC
Caprolactam	340	μg/kg	1800 U	890	J	NC	NC	400 U	380 U	NC	NC
Carbazole	660	μg/kg	800 J	480	J	NA	320	120 J	330 J	NA	210
Chrysene	660	μg/kg	2800	2100		28.6	NA	1200	4100 J	NA	2900
Dibenzo(a,h)anthracene	660	μg/kg	420 J	310		NA	110	180 J	650 J	NA	470
Dibenzofuran	340	μg/kg	190 J	120	J	NA	70	63 J	260 J	NA	197
Diethylphthalate	340	μg/kg	930 U	890	J	NC	NC	210 U	380 UJ	NC	NC
Dimethylphthalate	340	μg/kg	930 U	890	U	NC	NC	64 J	380 UJ	NA	316
Di-n-butylphthalate	340	μg/kg	930 U	890	J	NC	NC	210 U	380 UJ	NC	NC
Di-n-octylphthalate	340	μg/kg	1800 U	890	U	NC	NC	400 U	380 U	NC	NC
Fluoranthene	660	μg/kg	6200	4300		36.2	NA	2500	7500 J	100	NA
Fluorene	660	μg/kg	190 J	110	J	NA	80	140 J	360 J	NA	220
Hexachlorobenzene	340	μg/kg	930 U	250	UJ	NC	NC	210 U	100 UJ	NC	NC
Hexachlorobutadiene	340	μg/kg	930 U	330	UJ	NC	NC	210 U	150 UJ	NC	NC
Hexachlorocyclopentadiene	340	μg/kg	1800 U	2590	U	NC	NC	400 UJ	1100 U	NC	NC
Hexachloroethane	660	μg/kg	930 U	430	U	NC	NC	210 U	180 U	NC	NC
Indeno(1,2,3-cd)pyrene	340	μg/kg	1100 J	1300		16.7	NA	590	2800 J	NA	2210
Isophorone	340	μg/kg	930 U	300	UJ	NC	NC	210 U	120 UJ	NC	NC
Naphthalene	340	μg/kg	930 U	890	U	NC	NC	70 J	370 J	NA	300
Nitrobenzene	340	μg/kg	930 U	400		NC	NC	210 U	170 UJ	NC	NC
N-Nitroso-di-n-propylamine	340	μg/kg	930 U	230	U	NC	NC	210 U	90 UJ	NC	NC



Table 3-1a

Soil Samples Data Comparison - Metals, PCBs, and SVOCs Pierson's Creek Site, Operable Unit 2 Newark, New Jersey

	2 x		GS-B-112-001-CDN	Λ	GS-B-112-001-GEC	SY			GS-SS-101-001-CDM	Л G	GS-SS-101-001-GEOSY		
Chemical Name	CRQL	Unit	12/13/2018		12/13/2018		RPD	ABS	12/4/2018		12/4/2018	RPD	ABS
	CNQL		GS-B	3-11	12-001				GS-SS	S-10	01-001		
N-Nitrosodiphenylamine	340	μg/kg	930 U	J	210	UJ	NC	NC	210 U	J	90 UJ	NC	NC
Pentachlorophenol	660	μg/kg	1800 U	J	580	U	NC	NC	400 U	J	250 U	NC	NC
Phenanthrene	340	μg/kg	4700		2300	J	68.6	NA	1700		3800 J	76.4	NA
Phenol	660	μg/kg	1800 U	J	890	U	NC	NC	95 J		54 J	NA	41
Pyrene	340	μg/kg	5600		3500		46.2	NA	2700		7600 J	95.1	NA

Notes:

- 1. Sample pairs with RPD or ABS outside of criteria in the table are highlighted in yellow
- 2. PRP's laboratory analyzed 1,4-dioxane as a volatile organic compound.

ABS = absolute difference

CRQL = contract required quantitation limit

mg/kg = milligram per kilogram

NA = not applicable

NC = not calculable

PCB = polychlorinated biphenyl

RPD = relative percent difference

SVOC = semivolatile organic compound

μg/kg = microgram per kilogram

Data Validation Qualifiers

J = The analyte was positively identified, and the associated numerical value is the approximate concentration of the analyte in the sample.

J+ = The result is an estimated quantity, but the result may be biased high.

J- = The result is an estimated quantity, but the result may be biased low.

NJ = The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents its approximate concentration.

R = The sample results are unusable due to poor data quality; certain criteria were not met. The analyte may or may not be present in the sample at the reported concentration.

U = The analyte was not detected at a level greater than or equal to the level of the sample CRQL.



	2 x		GS-SS-101-002-C	M	GS-SS-101-002-GEC	SY			GS-SS-111-001-C	MC	GS-SS-111-001-G	EOSY		
Chemical Name	CRQL	Unit	12/4/2018		12/4/2018		RPD	ABS	12/4/2018		12/4/2018		RPD	ABS
	CKQL		GS	-SS-	101-002				GS	-SS-	111-001			
Metals														
Aluminum	40	mg/kg	6150		6160		0.16	NA	5390		4520		17.6	NA
Antimony	2	mg/kg	0.73	J	3.59	J	NA	2.86	0.92	J	2.99	J	NA	2.07
Arsenic	1	mg/kg	8.2		10.5		24.6	NA	13		22.1		51.9	NA
Barium	10	mg/kg	121		131		7.94	NA	90.1		78.4		13.9	NA
Beryllium	1	mg/kg	0.45		0.316	J	NA	0.13	0.42		0.279	J	NA	0.14
Cadmium	1	mg/kg	1.6		1.5		NA	0.1	0.59		0.643	J	NA	0.05
Calcium	1000	mg/kg	17800		39000		74.6	NA	7300		6480		11.9	NA
Chromium	2	mg/kg	46.3		31		39.6	NA	18.6		17.8		4.4	NA
Cobalt	1	mg/kg	25		20.6		19.3	NA	9.9		8.34		17.1	NA
Copper	2	mg/kg	85		74		13.8	NA	54.9		62.7		13.3	NA
Iron	20	mg/kg	18600		16200		13.8	NA	24100		26900		11	NA
Lead	1	mg/kg	613		449		30.9	NA	228		326		35.4	NA
Magnesium	1000	mg/kg	2430		2500		NA	70	2770		2390		NA	380
Manganese	1	mg/kg	289		272		6.06	NA	265		453		52.4	NA
Mercury	0	mg/kg	101		214		71.7	NA	5.4	J-	5.14		4.93	NA
Nickel	1	mg/kg	23.7		18.8		23.1	NA	14.6		13.3		9.32	NA
Potassium	1000	mg/kg	562		641		NA	79	448		443		NA	5
Selenium	5	mg/kg	3	כ	0.642	J	NA	2.36	2.8	UJ	0.719	J	NA	2.08
Silver	1	mg/kg	0.87	J	1.95		NA	1.08	0.8	IJ	0.516	J	NA	0.28
Sodium	1000	mg/kg	723		919		NA	196	1350		1430		NA	80
Thallium	1	mg/kg	0.8	J	1.86	J	NA	1.06	0.89	J	1.69	U	NA	0.8
Vanadium	5	mg/kg	26.7		23.8		11.5	NA	34.2		31		9.82	NA
Zinc	2	mg/kg	231		210		9.52	NA	190		766		121	NA
Polychlorinated Biphenyl														
Aroclor 1016	66	μg/kg	39	U	40	U	NC	NC	36	J	30	U	NC	NC
Aroclor 1221	66	μg/kg	39	U	40	U	NC	NC	36	U	30	U	NC	NC
Aroclor 1232	66	μg/kg	39	U	40	U	NC	NC	36	U	30	U	NC	NC
Aroclor 1242	66	μg/kg	39	U	73.3	JN	NA	34.3	36	U	30	U	NC	NC
Aroclor 1248	66	μg/kg	39	U	40	U	NC	NC	36	U	30	U	NC	NC
Aroclor 1254	66	μg/kg	160		417	J	NA	257	78	J	121	J	NA	43
Aroclor 1260	66	μg/kg	120	J	189	J	NA	69	85	NJ	83.5	J	NA	1.5
Aroclor 1262	66	μg/kg	39	U	40	U	NC	NC	36	U	30	U	NC	NC



	2 x		GS-SS-101-002-C	MC	GS-SS-101-002-GEC	SY			GS-SS-111-001-C	MC	GS-SS-111-001-G	EOSY		
Chemical Name	CRQL	Unit	12/4/2018		12/4/2018		RPD	ABS	12/4/2018		12/4/2018		RPD	ABS
	CKQL		GS	-SS-	101-002				GS	-SS-	-111-001			
Aroclor 1268	66	μg/kg	39	U	40	U	NC	NC	36	U	30	U	NC	NC
Semivolatile Organic Compound														
1,1'-Biphenyl	340	μg/kg	200	U	290	J	NA	90	190	U	204	J	NA	14
1,2,4,5-Tetrachlorobenzene	340	μg/kg	200	U	400	UJ	NC	NC	190	U	2090	UJ	NC	NC
1,4-Dioxane ²	340	μg/kg	79	UJ			NC	NC	73	UJ	890	UJ	NC	NC
2,2'-Oxybis(1-chloropropane)	340	μg/kg	390	U	470	UJ	NC	NC	360	U			NC	NC
2,3,4,6-Tetrachlorophenol	134	μg/kg	200	U	400	UJ	NC	NC	190	U	1100	UJ	NC	NC
2,4,5-Trichlorophenol	340	μg/kg	200	U	400	UJ	NC	NC	190	U	890	UJ	NC	NC
2,4,6-Trichlorophenol	340	μg/kg	200	U	230	UJ	NC	NC	190	U	890	UJ	NC	NC
2,4-Dichlorophenol	340	μg/kg	200	U	180	UJ	NC	NC	190	U	510	UJ	NC	NC
2,4-Dimethylphenol	340	μg/kg	200	U	370	UJ	NC	NC	190	U	430	UJ	NC	NC
2,4-Dinitrophenol	340	μg/kg	390	UJ	570	U	NC	NC	360	UJ	860	UJ	NC	NC
2,4-Dinitrotoluene	660	μg/kg	200	U	200	UJ	NC	NC	190	כ	1290	U	NC	NC
2,6-Dinitrotoluene	340	μg/kg	200	U	150	UJ	NC	NC	190	כ	440	UJ	NC	NC
2-Chloronaphthalene	340	μg/kg	200	U	400	IJ	NC	NC	190	J	360	UJ	NC	NC
2-Chlorophenol	340	μg/kg	200	U	120	UJ	NC	NC	190	J	890	UJ	NC	NC
2-Methylnaphthalene	340	μg/kg	100	J	790	J	NA	690	150	J	300	UJ	NA	150
2-Methylphenol	340	μg/kg	390	U	400	IJ	NC	NC	360	J	260	J	NA	100
2-Nitroaniline	660	μg/kg	200	U	400	UJ	NC	NC	190	J	890	UJ	NC	NC
2-Nitrophenol	340	μg/kg	200	U	870	UJ	NC	NC	190	J	890	UJ	NC	NC
3,3'-Dichlorobenzidine	340	μg/kg	390	U	310	UJ	NC	NC	360	U	2000	UJ	NC	NC
3-Nitroaniline	660	μg/kg	390	U	400	UJ	NC	NC	360	U	680	UJ	NC	NC
4,6-Dinitro-2-methylphenol	660	μg/kg	390	UJ	570	UJ	NC	NC	360	U	890		NC	NC
4-Bromophenyl-phenylether	660	μg/kg	200	U	400	UJ	NC	NC	190	U	1290	UJ	NC	NC
4-Chloro-3-methylphenol	340	μg/kg	200		400	UJ	NC	NC	190	U	890		NC	NC
4-Chloroaniline	340	μg/kg	390	U	400		NC	NC	360		890	UJ	NC	NC
4-Chlorophenyl-phenylether	660	μg/kg	200	U	400	UJ	NC	NC	190	U	890	UJ	NC	NC
4-Methylphenol	340	μg/kg	390				NC	NC	360	_	890	UJ	NC	NC
4-Nitroaniline	660	μg/kg	390	_	400	UJ	NC	NC	360	U			NC	NC
4-Nitrophenol	660	μg/kg	390	U	560	U	NC	NC	360	U	890	UJ	NC	NC
Acenaphthene	660	μg/kg	400		2200	_	NA	1800	190	_	1290	_	NC	NC
Acenaphthylene	340	μg/kg	200	U	480	J	NA	280	190	U	720	UJ	NC	NC



	2 x		GS-SS-101-002-C	M	GS-SS-101-002-GEOSY	,		GS-SS-111-001-CDM	GS-SS-111-001-GEOSY		
Chemical Name	CRQL	Unit	12/4/2018		12/4/2018	RPE	ABS	12/4/2018	12/4/2018	RPD	ABS
	CKQL		GS	-SS-	101-002			GS-SS	5-111-001		
Acetophenone	340	μg/kg	390	U	400 UJ	NC	NC	360 U	720 UJ	NC	NC
Anthracene	660	μg/kg	1000		5400 J	NA	4400	190 U	890 UJ	NC	NC
Atrazine	340	μg/kg	390	U	310 UJ	NC	NC	360 U	540 UJ	NC	NC
Benzaldehyde	660	μg/kg	390	U	520 U	NC	NC	360 U	720 UJ	NC	NC
Benzo(a)anthracene	660	μg/kg	1900		8600 J	128	NA	180 J	1200 U	NA	1020
Benzo(a)pyrene	340	μg/kg	1500		8300 J	139	NA	200	340 J	NA	140
Benzo(b)fluoranthene	340	μg/kg	1800		9800 J	138	NA	280	520 J	NA	240
Benzo(g,h,i)perylene	340	μg/kg	870		5400 J	144	NA	170 J	660 J	NA	490
Benzo(k)fluoranthene	340	μg/kg	600		2800 J	NA	2200	83 J	410 J	NA	327
Bis(2-chloroethoxy)methane	340	μg/kg	200	U	430 UJ	NC	NC	190 U	240 J	NA	50
Bis(2-chloroethyl)ether	340	μg/kg	390	U	140 UJ	NC	NC	360 U	980 UJ	NC	NC
Bis(2-ethylhexyl)phthalate	340	μg/kg	420		2900 J	NA	2480	190 U	310 UJ	NC	NC
Butylbenzylphthalate	660	μg/kg	200	U	400 UJ	NC	NC	190 U	890 UJ	NC	NC
Caprolactam	340	μg/kg	390	U	400 U	NC	NC	360 U	890 UJ	NC	NC
Carbazole	660	μg/kg	160	J	1000 J	NΑ	840	360 U	890 U	NC	NC
Chrysene	660	μg/kg	1800		7800 J	125	NA	230	890 UJ	NA	660
Dibenzo(a,h)anthracene	660	μg/kg	260		1300 J	NΑ	1040	190 U	400 J	NA	210
Dibenzofuran	340	μg/kg	260		1300 J	NA	1040	190 U	140 J	NA	50
Diethylphthalate	340	μg/kg	200	U	400 UJ	NC	NC	190 U	890 UJ	NC	NC
Dimethylphthalate	340	μg/kg	66	J	400 UJ	NA	334	89 J	890 UJ	NA	801
Di-n-butylphthalate	340	μg/kg	200	U	400 UJ	NC	NC	190 U	890 UJ	NC	NC
Di-n-octylphthalate	340	μg/kg	390	U	400 U	NC	NC	360 U	890 UJ	NC	NC
Fluoranthene	660	μg/kg	3300		14000 J	124	NA	300 J	890 U	NA	590
Fluorene	660	μg/kg	490		2800 J	NA	2310	190 U	470 J	NA	280
Hexachlorobenzene	340	μg/kg	200	U	110 UJ	NC	NC	190 U	890 UJ	NC	NC
Hexachlorobutadiene	340	μg/kg	200	U	150 UJ	NC	NC	190 U	250 UJ	NC	NC
Hexachlorocyclopentadiene	340	μg/kg	390	UJ	1200 U	NC	NC	360 UJ	340 UJ	NC	NC
Hexachloroethane	660	μg/kg	200	U	200 U	NC	NC	190 U	2590 U	NC	NC
Indeno(1,2,3-cd)pyrene	340	μg/kg	800		5700 J	NA	4900	130 J	430 U	NA	300
Isophorone	340	μg/kg	200	U	180 J	NΑ	20	190 U	400 J	NA	210
Naphthalene	340	μg/kg	130	J	1000 J	NΑ	870	190 U	890 UJ	NC	NC
Nitrobenzene	340	μg/kg	200	U	180 UJ	NC	NC	190 U	400 UJ	NC	NC
N-Nitroso-di-n-propylamine	340	μg/kg	200	U	100 UJ	NC	NC	190 U	230 UJ	NC	NC



Table 3-1a

Soil Samples Data Comparison - Metals, PCBs, and SVOCs Pierson's Creek Site, Operable Unit 2 Newark, New Jersey

	3,4	2 x Unit	GS-SS-101-002-CD	М	GS-SS-101-002-GEOSY			GS-SS-111-001-CDM	GS-SS-111-001-GEOSY		
Chemical Name	CRQL	Unit	12/4/2018		12/4/2018	RPD	ABS	12/4/2018	12/4/2018	RPD	ABS
	CKQL		GS-	SS-	101-002			GS-SS-	-111-001		
N-Nitrosodiphenylamine	340	μg/kg	200	С	90 UJ	NC	NC	190 U	210 UJ	NC	NC
Pentachlorophenol	660	μg/kg	390	U	270 U	NC	NC	360 U	600 U	NC	NC
Phenanthrene	340	μg/kg	3800		14000 J	115	NA	230	310 J	NA	80
Phenol	660	μg/kg	100	J	57 J	NA	43	100 J	890 U	NA	790
Pyrene	340	μg/kg	4800		15000 J	103	NA	380	460 J	NA	80

Notes:

- 1. Sample pairs with RPD or ABS outside of criteria in the table are highlighted in yellow
- 2. PRP's laboratory analyzed 1,4-dioxane as a volatile organic compound.

ABS = absolute difference

CRQL = contract required quantitation limit

mg/kg = milligram per kilogram

NA = not applicable

NC = not calculable

PCB = polychlorinated biphenyl

RPD = relative percent difference

SVOC = semivolatile organic compound

μg/kg = microgram per kilogram

Data Validation Qualifiers

J = The analyte was positively identified, and the associated numerical value is the approximate concentration of the analyte in the sample.

J+ = The result is an estimated quantity, but the result may be biased high.

J- = The result is an estimated quantity, but the result may be biased low.

NJ = The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents its approximate concentration.

R = The sample results are unusable due to poor data quality; certain criteria were not met. The analyte may or may not be present in the sample at the reported concentration.

U = The analyte was not detected at a level greater than or equal to the level of the sample CRQL.



Table 3-1a
Soil Samples Data Comparison - Metals, PCBs, and SVOCs
Pierson's Creek Site, Operable Unit 2
Newark, New Jersey

	2 11		GS-B-114-001-C	M	GS-B-114-001-GEO	SY		
Chemical Name	2 x	Unit	2/9/2019		2/9/2019		RPD	ABS
	CRQL		GS	-B-:	114-001			
Metals								
Aluminum	40	mg/kg	3940		3950		0.25	NA
Antimony	2	mg/kg	5.7	J	31.2		138	NA
Arsenic	1	mg/kg	230		254		9.92	NA
Barium	10	mg/kg	260		144		57.4	NA
Beryllium	1	mg/kg	0.37	J	0.192	J	NA	0.18
Cadmium	1	mg/kg	46.4		22.1		70.9	NA
Calcium	1000	mg/kg	5680		5670		0.18	NA
Chromium	2	mg/kg	274		316		14.2	NA
Cobalt	1	mg/kg	11.2		8.84		23.6	NA
Copper	2	mg/kg	1400		1380		1.44	NA
Iron	20	mg/kg	20000		14700		30.5	NA
Lead	1	mg/kg	1360		1200		12.5	NA
Magnesium	1000	mg/kg	1260		1440		NA	180
Manganese	1	mg/kg	121		95.5		23.6	NA
Mercury	0	mg/kg	128		97.9		26.6	NA
Nickel	1	mg/kg	67.7		64.1		5.46	NA
Potassium	1000	mg/kg	636	כ	552		NA	84
Selenium	5	mg/kg	4.5	UJ	2.08	J	NA	2.42
Silver	1	mg/kg	61		75.4		21.1	NA
Sodium	1000	mg/kg	636	J	319		NA	317
Thallium	1	mg/kg	1.4	J	2.74	U	NA	1.34
Vanadium	5	mg/kg	38.3		58.2		41.2	NA
Zinc	2	mg/kg	1510		1010		39.7	NA
Polychlorinated Biphenyl								
Aroclor 1016	66	μg/kg	52000		28390	U	58.7	NA
Aroclor 1221	66	μg/kg	560	U	28390	U	NC	NC
Aroclor 1232	66	μg/kg	560	U	28390	U	NC	NC
Aroclor 1242	66	μg/kg	560	U	348000	J	199	NA
Aroclor 1248	66	μg/kg	560	U	28390	U	NC	NC
Aroclor 1254	66	μg/kg	560	U	28390	U	NC	NC
Aroclor 1260	66	μg/kg	34000		131000	J	118	NA
Aroclor 1262	66	μg/kg	560	U	28390	U	NC	NC



Table 3-1a
Soil Samples Data Comparison - Metals, PCBs, and SVOCs
Pierson's Creek Site, Operable Unit 2
Newark, New Jersey

	2 x		GS-B-114-001-C	MC	GS-B-114-001-GE0	OSY		
Chemical Name	CRQL	Unit	2/9/2019		2/9/2019		RPD	ABS
	CKQL		GS	-B-:	114-001			
Aroclor 1268	66	μg/kg	560	U	28390	U	NC	NC
Semivolatile Organic Compound								
1,1'-Biphenyl	340	μg/kg	1400	U	280	J	NA	1120
1,2,4,5-Tetrachlorobenzene	340	μg/kg	1400	U	540	J	NA	860
1,4-Dioxane ²	340	μg/kg	570	R	1400000	R	NC	NC
2,2'-Oxybis(1-chloropropane)	340	μg/kg	2800	U	1390	U	NC	NC
2,3,4,6-Tetrachlorophenol	134	μg/kg	1400	U	1100	U	NC	NC
2,4,5-Trichlorophenol	340	μg/kg	1400	٦	1100	U	NC	NC
2,4,6-Trichlorophenol	340	μg/kg	1400	U	640	U	NC	NC
2,4-Dichlorophenol	340	μg/kg	1400	U	550	U	NC	NC
2,4-Dimethylphenol	340	μg/kg	1400	J	1100	U	NC	NC
2,4-Dinitrophenol	340	μg/kg	2800	J	1600	U	NC	NC
2,4-Dinitrotoluene	660	μg/kg	1400	٦	560	U	NC	NC
2,6-Dinitrotoluene	340	μg/kg	1400	IJ	440	U	NC	NC
2-Chloronaphthalene	340	μg/kg	1400	J	1100	U	NC	NC
2-Chlorophenol	340	μg/kg	1400	J	370	U	NC	NC
2-Methylnaphthalene	340	μg/kg	570	J	1600		NA	1030
2-Methylphenol	340	μg/kg	2800	J	1100	U	NC	NC
2-Nitroaniline	660	μg/kg	1400	U	1100	U	NC	NC
2-Nitrophenol	340	μg/kg	1400	J	2500	U	NC	NC
3,3'-Dichlorobenzidine	340	μg/kg	2800	J	870	UJ	NC	NC
3-Nitroaniline	660	μg/kg	2800	UJ	1100	UJ	NC	NC
4,6-Dinitro-2-methylphenol	660	μg/kg	2800	U	1600		NC	NC
4-Bromophenyl-phenylether	660	μg/kg	1400	U	1100		NC	NC
4-Chloro-3-methylphenol	340	μg/kg	1400	U	1100		NC	NC
4-Chloroaniline	340	μg/kg	2800	J	1100	U	NC	NC
4-Chlorophenyl-phenylether	660	μg/kg	1400	U	1100	U	NC	NC
4-Methylphenol	340	μg/kg	2800				NC	NC
4-Nitroaniline	660	μg/kg	2800	U	1100		NC	NC
4-Nitrophenol	660	μg/kg	2800	U	1600	U	NC	NC
Acenaphthene	660	μg/kg	1400	U	500	J	NA	900
Acenaphthylene	340	μg/kg	1400	U	280	J	NA	1120



Table 3-1a
Soil Samples Data Comparison - Metals, PCBs, and SVOCs
Pierson's Creek Site, Operable Unit 2
Newark, New Jersey

Chemical Name	2 x	Unit	GS-B-114-001-CD 2/9/2019	M	GS-B-114-001-GEO	SY	RPD	ABS
	CRQL			-B-:	114-001			
Acetophenone	340	μg/kg	2800	U	1100	U	NC	NC
Anthracene	660	μg/kg	14000		40000		96.3	NA
Atrazine	340	μg/kg	2800	U	910 (U	NC	NC
Benzaldehyde	660	μg/kg	2800	U	1500 (U	NC	NC
Benzo(a)anthracene	660	μg/kg	970	J	2100		NA	1130
Benzo(a)pyrene	340	μg/kg	610	J	1400		NA	790
Benzo(b)fluoranthene	340	μg/kg	1200	J	2700		76.9	NA
Benzo(g,h,i)perylene	340	μg/kg	470	J	1300		NA	830
Benzo(k)fluoranthene	340	μg/kg	400	J	870		NA	470
Bis(2-chloroethoxy)methane	340	μg/kg	1400	U	1200 (U	NC	NC
Bis(2-chloroethyl)ether	340	μg/kg	2800	U	400 (U	NC	NC
Bis(2-ethylhexyl)phthalate	340	μg/kg	9000		17000		61.5	NA
Butylbenzylphthalate	660	μg/kg	1400	U	1100 (U	NC	NC
Caprolactam	340	μg/kg	2800	U	1100 (U	NC	NC
Carbazole	660	μg/kg	2800	U	420 J	J	NA	2380
Chrysene	660	μg/kg	1200	J	2400		NA	1200
Dibenzo(a,h)anthracene	660	μg/kg	1400	כ	450		NA	950
Dibenzofuran	340	μg/kg	1400	כ	410 J	J	NA	990
Diethylphthalate	340	μg/kg	1400	כ	1100 (U	NC	NC
Dimethylphthalate	340	μg/kg	1400	כ	1100	U	NC	NC
Di-n-butylphthalate	340	μg/kg	580	J	1600		NA	1020
Di-n-octylphthalate	340	μg/kg	680	J	1900		NA	1220
Fluoranthene	660	μg/kg	2300	J	5000		74	NA
Fluorene	660	μg/kg	340	J	790 J	J	NA	450
Hexachlorobenzene	340	μg/kg	1400	J	330 (U	NC	NC
Hexachlorobutadiene	340	μg/kg	1400	כ	410 (UJ	NC	NC
Hexachlorocyclopentadiene	340	μg/kg	2800	J	3290 เ	U	NC	NC
Hexachloroethane	660	μg/kg	1400	J	550 l	U	NC	NC
Indeno(1,2,3-cd)pyrene	340	μg/kg	440	J	1300		NA	860
Isophorone	340	μg/kg	1400	U	370 l	U	NC	NC
Naphthalene	340	μg/kg	3100		8800		95.8	NA
Nitrobenzene	340	μg/kg	1400	U	500 l	U	NC	NC
N-Nitroso-di-n-propylamine	340	μg/kg	1400	U	280 l	U	NC	NC



Table 3-1a

Soil Samples Data Comparison - Metals, PCBs, and SVOCs Pierson's Creek Site, Operable Unit 2 Newark, New Jersey

	2 x		GS-B-114-001-C	MC	GS-B-114-001-GEO	SY		
Chemical Name	CRQL	Unit	2/9/2019		2/9/2019		RPD	ABS
	CKQL		GS	-B-:	114-001			
N-Nitrosodiphenylamine	340	μg/kg	1400	٦	280 l	J	NC	NC
Pentachlorophenol	660	μg/kg	2800	U	750 l	J	NC	NC
Phenanthrene	340	μg/kg	1600		3700		79.2	NA
Phenol	660	μg/kg	2800	U	1100 l	J	NC	NC
Pyrene	340	μg/kg	1800		4200		80	NA

Notes:

- 1. Sample pairs with RPD or ABS outside of criteria in the table are highlighted in yellow
- 2. PRP's laboratory analyzed 1,4-dioxane as a volatile organic compound.

ABS = absolute difference

CRQL = contract required quantitation limit

mg/kg = milligram per kilogram

NA = not applicable

NC = not calculable

PCB = polychlorinated biphenyl

RPD = relative percent difference

SVOC = semivolatile organic compound

μg/kg = microgram per kilogram

Data Validation Qualifiers

J = The analyte was positively identified, and the associated numerical value is the approximate concentration of the analyte in the sample.

J+ = The result is an estimated quantity, but the result may be biased high.

J- = The result is an estimated quantity, but the result may be biased low.

NJ = The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents its approximate concentration.

R = The sample results are unusable due to poor data quality; certain criteria were not met. The analyte may or may not be present in the sample at the reported concentration.

U = The analyte was not detected at a level greater than or equal to the level of the sample CRQL.



	2 x		GS-B-101-004-CDM	GS-B-101-004-GEOSY			GS-B-102-004-CDM	GS-B-102-004-GEOS	Υ	
Chemical Name	CRQL	Unit	12/14/2018	12/14/2018	RPD	ABS	12/17/2018	12/17/2018	RPE	ABS
	CKQL		GS-B-1	01-004			GS-B-2	102-004		
Volatile Organic Compound										
1,1,1-Trichloroethane	10	μg/kg	5 U	20 U	NC	NC	12 U	90 U	NC.	. NC
1,1,2,2-Tetrachloroethane	10	μg/kg	5 UJ	20 U	NC	NC	12 U	90 U	NC.	NC NC
1,1,2-Trichloro-1,2,2-trifluoroethane	10	μg/kg	5 U	230 U	NC	NC	12 U	730 U	NC	NC NC
1,1,2-Trichloroethane	10	μg/kg	5 U	50 U	NC	NC	12 U	180 U	NC	NC NC
1,1-Dichloroethane	10	μg/kg	5 UJ	50 U	NC	NC	12 U	180 U	NC	NC NC
1,1-Dichloroethene	10	μg/kg	5 U	50 U	NC	NC	12 U	180 U	NC	NC NC
1,2,3-Trichlorobenzene	10	μg/kg	5 UJ	110 U	NC	NC	12 UJ	360 U	NC	NC NC
1,2,4-Trichlorobenzene	10	μg/kg	5 UJ	110 U	NC	NC	12 UJ	64 J	NA	52
1,2-Dibromo-3-chloropropane	10	μg/kg	5 UJ	180 U	NC	NC	12 UJ	540 U	NC	. NC
1,2-Dibromoethane	10	μg/kg	5 U	50 U	NC	NC	12 U	180 U	NC	NC NC
1,2-Dichlorobenzene	10	μg/kg	4 J+	13 J	NA	9	12 UJ	360 U	NC	. NC
1,2-Dichloroethane	10	μg/kg	5 U	50 U	NC	NC	12 U	180 U	NC.	NC NC
1,2-Dichloropropane	10	μg/kg	5 U	50 U	NC	NC	12 U	180 U	NC	NC NC
1,3-Dichlorobenzene	10	μg/kg	5 UJ	110 U	NC	NC	12 UJ	360 U	NC.	NC NC
1,4-Dichlorobenzene	10	μg/kg	2 J+	15 J	NA	13	12 UJ	80 J	NA	68
1,4-Dioxane ²	10	μg/kg		5900 R	NC	NC		18000 R	NC	NC NC
2-Butanone	20	μg/kg	6.6 J	580 U	NA	573	96	1790 U	180	NA
2-Hexanone	20	μg/kg	9.9 U	580 U	NC	NC	24 U	1790 U	NC	NC NC
4-Methyl-2-pentanone	20	μg/kg	9.9 U	580 U	NC	NC	24 U	1790 U	NC	. NC
Acetone	20	μg/kg	65	580 U	160	NA	290	1790 U	144	l NA
Benzene	10	μg/kg	210 J	1800	158	NA	31	580	180) NA
Bromochloromethane	10	μg/kg	5 UJ	110 U	NC	NC	12 U	360 U	NC	NC NC
Bromodichloromethane	10	μg/kg	5 U	20 U	NC	NC	12 U	90 U	NC.	NC NC
Bromoform	10	μg/kg	5 UJ	230 U	NC	NC	12 UJ	730 U	NC.	NC NC
Bromomethane	10	μg/kg	5 U	110 U	NC	NC	12 U	360 U	NC	. NC
Carbon Disulfide	10	μg/kg	11	580 U	NA	569	12	1790 U	N/A	<u>1778</u>
Carbon Tetrachloride	10	μg/kg	5 U	50 U	NC	NC	12 U	180 U	NC	. NC
Chlorobenzene	10	μg/kg	4.3 J	40	NA	35.7	12 U	160	NA	148
Chloroethane	10	μg/kg	5 U	110 U	NC	NC	12 U	360 U		
Chloroform	10	μg/kg	42 J-	80 U	62.3	NA	12 U	270 U	NC.	: NC
Chloromethane	10	μg/kg	5 U	230 U	NC	NC	12 U	730 U	NC.	: NC
cis-1,2-Dichloroethene	10	μg/kg	5 U	50 U	NC	NC	12 U	54 J	NA	42
cis-1,3-Dichloropropene	10	μg/kg	5 U	20 U	NC	NC	12 U	90 U	NC.	: NC
Cyclohexane	10	μg/kg	12 J+	69 J	NA	57	12 U	1790 U	NC.	: NC
Dibromochloromethane	10	μg/kg	5 UJ	50 U	NC	NC	12 U	180 U	NC.	NC NC



	2 x		GS-B-101-004-CDM	GS-B-101-004-GEOSY	′		GS-B-102-004-CD	М	GS-B-102-004-GEOS	Υ	
Chemical Name	CRQL	Unit	12/14/2018	12/14/2018	RPD	ABS	12/17/2018		12/17/2018	RPD	ABS
	CKQL		GS-B-	101-004			GS-	B-1	.02-004		
Dichlorodifluoromethane	10	μg/kg	5 U	580 U	NC	NC	12	U	1790 U	NC	NC
Ethylbenzene	10	μg/kg	16	100	NA	84	4.9	J	150 J	NA	145
Isopropylbenzene	10	μg/kg	20	54 J	NA	34	9.7	J	200	NA	190
m,p-Xylene	10	μg/kg	8.7	74 J	NA	65.3	13		330 J	NA	317
Methyl acetate	10	μg/kg	5 U	1200	NA	1195	12	U	280 J	NA	268
Methyl tert-Butyl Ether	10	μg/kg	5 U	110 U	NC	NC	12	U	360 U	NC	NC
Methylcyclohexane	10	μg/kg	1.1 J+	230 U	NA	229	5.9	J+	120 J	NA	114
Methylene Chloride	10	μg/kg	10	280 U	NA	270	12	U	910 U	NC	NC
o-Xylene	10	μg/kg	24	69	NA	45	11	J	220	NA	209
Styrene	10	μg/kg	5 U	50 U	NC	NC	12	U	180 U	NC	NC
Tetrachloroethene	10	μg/kg	0.96 J	20 U	NA	19	12	U	90 U	NC	NC
Toluene	10	μg/kg	3 J	50 U	NA	47	9.9	J	230	NA	220
trans-1,2-Dichloroethene	10	μg/kg	5 U	80 U	NC	NC	12	U	270 U	NC	NC
trans-1,3-Dichloropropene	10	μg/kg	5 U	50 U	NC	NC	12	U	180 U	NC	NC
Trichloroethene	10	μg/kg	3.3 J	20 U	NA	16.7	12	U	110	NA	98
Trichlorofluoromethane	10	μg/kg	5 U	230 U	NC	NC	12	U	730 U	NC	NC
Vinyl Chloride	10	μg/kg	5 U	50 U	NC	NC	12	U	180 U	NC	NC

Notes:

- 1. Sample pairs with RPD or ABS outside of criteria in the table are highlighted in yellow
- 2. CDM Smith's assigned laboratory analyzed
- 1,4-dioxane as a semivolatile organic compound.

ABS = absolute difference

CRQL = contract required quantitation limit

mg/kg = milligram per kilogram

NA = not applicable

NC = not calculable

RPD = relative percent difference

μg/kg = microgram per kilogram

VOC = volatile organic compound

Data Validation Qualifiers

J = The analyte was positively identified, and the associated numerical value is the approximate concentration of the analyte in the sample.

J+ = The result is an estimated quantity, but the result may be biased high.

J- = The result is an estimated quantity, but the result may be biased low.

NJ = The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents its approximate concentration.

R = The sample results are unusable due to poor data quality; certain criteria were not met. The analyte may or may not be present in the sample at the reported concentration.

 $\mbox{\bf U}$ = The analyte was not detected at a level greater than or equal to the level of the sample CRQL.



	2		GS-B-106-004-CDM	GS-B-106-004-GEOSY			GS-B-111-004-CDM	1 GS-B-111-004-GEOSY	<i>'</i>	
Chemical Name	2 x CRQL	Unit	12/13/2018	12/13/2018	RPD	ABS	12/17/2018	12/17/2018	RPD	ABS
	CKQL		GS-B-1	06-004			GS-B	-111-004		
Volatile Organic Compound										
1,1,1-Trichloroethane	10	μg/kg	470 U	2590 U	NC	NC	400 L	2000 U	NC	NC
1,1,2,2-Tetrachloroethane	10	μg/kg	470 U	2590 U	NC	NC	400 L	2000 U	NC	NC
1,1,2-Trichloro-1,2,2-trifluoroethane	10	μg/kg	470 U	21000 U	NC	NC	400 L	16000 U	NC	NC
1,1,2-Trichloroethane	10	μg/kg	470 U	5190 U	NC	NC	400 L	4090 U	NC	NC
1,1-Dichloroethane	10	μg/kg	470 U	5190 U	NC	NC	400 L	4090 U	NC	NC
1,1-Dichloroethene	10	μg/kg	470 U	5190 U	NC	NC	400 L	4090 U	NC	NC
1,2,3-Trichlorobenzene	10	μg/kg	470 U	10000 U	NC	NC	400 L	8190 U	NC	NC
1,2,4-Trichlorobenzene	10	μg/kg	470 U	10000 U	NC	NC	400 L	8190 U	NC	NC
1,2-Dibromo-3-chloropropane	10	μg/kg	470 U	16000 U	NC	NC	400 L	12000 U	NC	NC
1,2-Dibromoethane	10	μg/kg	470 U	5190 U	NC	NC	400 L	4090 U	NC	NC
1,2-Dichlorobenzene	10	μg/kg	470 U	10000 U	NC	NC	400 L	8190 U	NC	NC
1,2-Dichloroethane	10	μg/kg	470 U	5190 U	NC	NC	400 L	4090 U	NC	NC
1,2-Dichloropropane	10	μg/kg	470 U	5190 U	NC	NC	400 L	4090 U	NC	NC
1,3-Dichlorobenzene	10	μg/kg	470 U	10000 U	NC	NC	400 L	8190 U	NC	NC
1,4-Dichlorobenzene	10	μg/kg	470 U	10000 U	NC	NC	400 L	8190 U	NC	NC
1,4-Dioxane ²	10	μg/kg		520000 R	NC	NC		410000 R	NC	NC
2-Butanone	20	μg/kg	940 U	52000 U	NC	NC	800 L	41000 U	NC	NC
2-Hexanone	20	μg/kg	940 U	52000 U	NC	NC	800 L	J 41000 U	NC	NC
4-Methyl-2-pentanone	20	μg/kg	940 U	52000 U	NC	NC	800 L	41000 U	NC	NC
Acetone	20	μg/kg	940 U	52000 U	NC	NC	800 L	41000 U	NC	NC
Benzene	10	μg/kg	320 J	6100	180	NA	160 J	2000 U	170	NA
Bromochloromethane	10	μg/kg	470 U	10000 U	NC	NC	400 L	8190 U	NC	NC
Bromodichloromethane	10	μg/kg	470 U	2590 U	NC	NC	400 L	2000 U	NC	NC
Bromoform	10	μg/kg	470 U	21000 U	NC	NC	400 L	16000 U	NC	NC
Bromomethane	10	μg/kg	470 U	10000 U	NC	NC	400 L	8190 U	NC	NC
Carbon Disulfide	10	μg/kg	470 U	52000 U	NC	NC	400 L	41000 U	NC	NC
Carbon Tetrachloride	10	μg/kg	470 U	5190 U	NC	NC	400 L	J 4090 U	NC	NC
Chlorobenzene	10	μg/kg	120 J	1100 J	161	NA	400 L	2000 U	NC	NC
Chloroethane	10	μg/kg	470 U	10000 U	NC	NC	400 L	8190 U	NC	NC
Chloroform	10	μg/kg	470 U	7800 U	NC	NC	400 L	6090 U	NC	NC
Chloromethane	10	μg/kg	470 U	21000 U	NC	NC	400 L	16000 U	NC	NC
cis-1,2-Dichloroethene	10	μg/kg	470 U	5190 U	NC	NC	400 L	4090 U	NC	NC
cis-1,3-Dichloropropene	10	μg/kg	470 U	2590 U	NC	NC	400 L	2000 U	NC	NC
Cyclohexane	10	μg/kg	470 U	52000 U	NC	NC	400 L	41000 U	NC	NC
Dibromochloromethane	10	μg/kg	470 U	5190 U	NC	NC	400 L	4090 U	NC	NC



	2 x		GS-B-106-004-CDM	GS-B-106-004-GEOSY			GS-B-111-004-CDN	M	GS-B-111-004-GEO	SY		
Chemical Name	CRQL	Unit	12/13/2018	12/13/2018	RPD	ABS	12/17/2018		12/17/2018		RPD	ABS
	CKQL		GS-B-1	06-004			GS-B	3-1:	11-004			
Dichlorodifluoromethane	10	μg/kg	470 U	52000 U	NC	NC	400	U	41000 L	J	NC	NC
Ethylbenzene	10	μg/kg	220 J	2000 J	160	NA	470000		380000		21.2	NA
Isopropylbenzene	10	μg/kg	10000	76000	153	NA	34000		41000		18.7	NA
m,p-Xylene	10	μg/kg	320 J	10000 U	188	NA	3100		7100 J		78.4	NA
Methyl acetate	10	μg/kg	470 U	21000 U	NC	NC	400	U	16000 L	J	NC	NC
Methyl tert-Butyl Ether	10	μg/kg	470 U	10000 U	NC	NC	400	U	8190 L	J	NC	NC
Methylcyclohexane	10	μg/kg	470 U	21000 U	NC	NC	400	U	16000 L	J	NC	NC
Methylene Chloride	10	μg/kg	470 U	26000 U	NC	NC	400	U	20000 L	J	NC	NC
o-Xylene	10	μg/kg	1600	13000	156	NA	13000		31000		81.8	NA
Styrene	10	μg/kg	470 U	5190 U	NC	NC	400	U	4090 L	J	NC	NC
Tetrachloroethene	10	μg/kg	470 U	2590 U	NC	NC	400	U	2000 L	J	NC	NC
Toluene	10	μg/kg	470 U	5190 U	NC	NC	400	U	4090 L	J	NC	NC
trans-1,2-Dichloroethene	10	μg/kg	470 U	7800 U	NC	NC	400	U	6090 L	J	NC	NC
trans-1,3-Dichloropropene	10	μg/kg	470 U	5190 U	NC	NC	400	U	4090 L	J	NC	NC
Trichloroethene	10	μg/kg	470 U	2590 U	NC	NC	400	U	2000 L	J	NC	NC
Trichlorofluoromethane	10	μg/kg	470 U	21000 U	NC	NC	400	U	16000 L	J	NC	NC
Vinyl Chloride	10	μg/kg	470 U	5190 U	NC	NC	400	U	4090 L	J	NC	NC

Notes:

- 1. Sample pairs with RPD or ABS outside of criteria in the table are highlighted in yellow
- 2. CDM Smith's assigned laboratory analyzed
 - 1,4-dioxane as a semivolatile organic compound.

ABS = absolute difference

CRQL = contract required quantitation limit

mg/kg = milligram per kilogram

NA = not applicable

NC = not calculable

RPD = relative percent difference

μg/kg = microgram per kilogram

VOC = volatile organic compound

Data Validation Qualifiers

J = The analyte was positively identified, and the associated numerical value is the approximate concentration of the analyte in the sample.

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R = The sample results are unusable due to poor data quality; certain criteria were not met. The analyte may or may not be present in the sample at the reported concentration.

U = The analyte was not detected at a level greater than or equal to the level of the sample CRQL.



Table 3-1b Soil Samples Data Comparison - VOCs Pierson's Creek Site, Operable Unit 2 Newark, New Jersey

	2		GS-B-112-004-CDM	GS-B-112-004-GEOSY			GS-SS-101-003-CD	М	GS-SS-101-003-GEOSY		T
Chemical Name	2 x CRQL	Unit	12/13/2018	12/13/2018	RPD	ABS	12/4/2018		12/4/2018	RPD	ABS
	CKQL		GS-B-1	112-004			GS-	SS-	101-003		
Volatile Organic Compound											
1,1,1-Trichloroethane	10	μg/kg	6 U	0.84 U	NC	NC	4.7	U	0.58 U	NC	NC
1,1,2,2-Tetrachloroethane	10	μg/kg	6 U	0.84 U	NC	NC	4.7	U	0.58 U	NC	NC
1,1,2-Trichloro-1,2,2-trifluoroethane	10	μg/kg	6 U	6.80 U	NC	NC	4.7	U	4.70 U	NC	NC
1,1,2-Trichloroethane	10	μg/kg	6 U	1.70 UJ	NC	NC	4.7	U	1.20 U	NC	NC
1,1-Dichloroethane	10	μg/kg	6 U	1.70 U	NC	NC	4.7	U	1.20 U	NC	NC
1,1-Dichloroethene	10	μg/kg	6 U	1.70 U	NC	NC	4.7	U	1.20 U	NC	NC
1,2,3-Trichlorobenzene	10	μg/kg	6 UJ	3.40 U	NC	NC	4.7	U	2.30 U	NC	NC
1,2,4-Trichlorobenzene	10	μg/kg	6 UJ	3.40 U	NC	NC	4.7	U	2.30 U	NC	NC
1,2-Dibromo-3-chloropropane	10	μg/kg	6 UJ	5.10 U	NC	NC	4.7	U	3.50 U	NC	NC
1,2-Dibromoethane	10	μg/kg	6 U	1.70 UJ	NC	NC	4.7	\supset	1.20 U	NC	NC
1,2-Dichlorobenzene	10	μg/kg	6 UJ	0.62 J	NA	5.38	4.7	\supset	2.30 U	NC	NC
1,2-Dichloroethane	10	μg/kg	6 U	1.70 U	NC	NC	4.7	\supset	1.20 U	NC	NC
1,2-Dichloropropane	10	μg/kg	6 U	1.70 U	NC	NC	4.7	\supset	1.20 U	NC	NC
1,3-Dichlorobenzene	10	μg/kg	6 UJ	3.40 U	NC	NC	4.7	\supset	2.30 U	NC	NC
1,4-Dichlorobenzene	10	μg/kg	6 UJ	3.40 U	NC	NC	4.7	\supset	2.30 U	NC	NC
1,4-Dioxane ²	10	μg/kg		170 R	NC	NC			110 R	NC	NC
2-Butanone	20	μg/kg	12 U	18	NA	6	9.4	U	10 U	NC	NC
2-Hexanone	20	μg/kg	12 U	10 UJ	NC	NC	9.4	U	10 U	NC	NC
4-Methyl-2-pentanone	20	μg/kg	12 U	10 UJ	NC	NC	9.4	U	10 U	NC	NC
Acetone	20	μg/kg	39	97	NA	58	24		17	NA	7
Benzene	10	μg/kg	6 U	0.59 J	NA	5.41	4.7	\supset	1.2	NA	3.5
Bromochloromethane	10	μg/kg	6 U	3.40 U	NC	NC	4.7	\supset	2.30 U	NC	NC
Bromodichloromethane	10	μg/kg	6 U	0.84 U	NC	NC	4.7	\supset	0.58 U	NC	NC
Bromoform	10	μg/kg	6 UJ	6.80 U	NC	NC	4.7	\supset	4.70 U	NC	NC
Bromomethane	10	μg/kg	6 U	3.40 U	NC	NC	4.7	U	2.30 U	NC	NC
Carbon Disulfide	10	μg/kg	6 U	10 U	NC	NC	1.8	J	10 U	NA	8.2
Carbon Tetrachloride	10	μg/kg	6 U	1.70 U	NC	NC	4.7	U	1.200 U	NC	NC
Chlorobenzene	10	μg/kg	6 U	0.84 UJ	NC	NC	4.7	J	0.580 U	NC	NC
Chloroethane	10	μg/kg	6 U	3.40 U	NC	NC	4.7	\supset	2.300 U	NC	NC
Chloroform	10	μg/kg	6 U	2.50 U	NC	NC	4.7	\supset	1.700 U	NC	NC
Chloromethane	10	μg/kg	6 U	6.80 U	NC	NC	4.7	U	4.700 U	NC	NC
cis-1,2-Dichloroethene	10	μg/kg	6 U	0.6 J	NA	5.4	4.7	U	1.200 U	NC	NC
cis-1,3-Dichloropropene	10	μg/kg	6 U	0.84 U	NC	NC	4.7	U	0.580 U	NC	NC
Cyclohexane	10	μg/kg	2.3 J	5.1 J	NA	2.8	4.7	U	10 U	NC	NC
Dibromochloromethane	10	μg/kg	6 U	1.70 UJ	NC	NC	4.7	U	1.20 U	NC	NC



	2 x		GS-B-112-004-CDM	GS-B-112-004-GEOSY			GS-SS-101-003-CD	М	GS-SS-101-003-GEOSY		
Chemical Name	CRQL	Unit	12/13/2018	12/13/2018	RPD	ABS	12/4/2018		12/4/2018	RPD	ABS
	CKQL		GS-B-1	12-004			GS-	SS-	101-003		
Dichlorodifluoromethane	10	μg/kg	6 U	10 U	NC	NC	4.7	U	10 U	NC	NC
Ethylbenzene	10	μg/kg	6 U	1.70 UJ	NC	NC	4.7	U	0.17 J	NA	4.53
Isopropylbenzene	10	μg/kg	6 U	2	NA	4	0.86	J	0.16 J	NA	0.7
m,p-Xylene	10	μg/kg	6 U	3.40 UJ	NC	NC	4.7	U	2.30 U	NC	NC
Methyl acetate	10	μg/kg	6 U	6.80 U	NC	NC	4.7	U	4.70 U	NC	NC
Methyl tert-Butyl Ether	10	μg/kg	6 U	0.39 J	NA	5.61	4.7	U	2.30 U	NC	NC
Methylcyclohexane	10	μg/kg	6 U	3.2 J	NA	2.8	4.7	U	4.70 U	NC	NC
Methylene Chloride	10	μg/kg	6 U	8.40 U	NC	NC	3	J	5.80 U	NA	2.8
o-Xylene	10	μg/kg	6 U	0.62 J	NA	5.38	4.7	U	1.20 U	NC	NC
Styrene	10	μg/kg	6 U	1.70 UJ	NC	NC	4.7	U	1.20 U	NC	NC
Tetrachloroethene	10	μg/kg	6 U	0.84 UJ	NC	NC	4.7	U	0.32 J	NA	4.38
Toluene	10	μg/kg	6 U	1.70 UJ	NC	NC	4.7	U	1.20 U	NC	NC
trans-1,2-Dichloroethene	10	μg/kg	6 U	2.50 U	NC	NC	4.7	U	1.70 U	NC	NC
trans-1,3-Dichloropropene	10	μg/kg	6 U	1.70 UJ	NC	NC	4.7	U	1.20 U	NC	NC
Trichloroethene	10	μg/kg	6 U	0.84 U	NC	NC	4.7	U	0.39 J	NA	4.31
Trichlorofluoromethane	10	μg/kg	6 U	6.80 U	NC	NC	4.7	U	4.70 U	NC	NC
Vinyl Chloride	10	μg/kg	6 U	1.70 U	NC	NC	4.7	U	1.20 U	NC	NC

Notes:

- 1. Sample pairs with RPD or ABS outside of criteria in the table are highlighted in yellow
- 2. CDM Smith's assigned laboratory analyzed 1,4-dioxane as a semivolatile organic compound.

ABS = absolute difference

CRQL = contract required quantitation limit

mg/kg = milligram per kilogram

NA = not applicable

NC = not calculable

RPD = relative percent difference

μg/kg = microgram per kilogram

VOC = volatile organic compound

Data Validation Qualifiers

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J+ = The result is an estimated quantity, but the result may be biased high.

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R = The sample results are unusable due to poor data quality; certain criteria were not met. The analyte may or may not be present in the sample at the reported concentration.

 $\mbox{\bf U}$ = The analyte was not detected at a level greater than or equal to the level of the sample CRQL.



	2		GS-SS-111-003-CD	М	GS-SS-111-003-G	EOSY			GS-B-114-001-CD	M	GS-B-114-001-GEO	SY		,
Chemical Name	2 x CRQL	Unit	12/4/2018		12/4/2018		RPD	ABS	2/9/2019		2/9/2019		RPD	ABS
	CKQL		GS-	SS-	-111-003				GS-	-B-:	114-001			
Volatile Organic Compound														
1,1,1-Trichloroethane	10	μg/kg	270	U	50.00	U	NC	NC	7400	J	7200		2.74	NA
1,1,2,2-Tetrachloroethane	10	μg/kg	270	U	50.00	U	NC	NC	12000	U	6800 l	U	NC	NC
1,1,2-Trichloro-1,2,2-trifluoroethane	10	μg/kg	270	U	400.00	U	NC	NC	12000	U	54000 l	U	NC	NC
1,1,2-Trichloroethane	10	μg/kg	270	U	100.00	U	NC	NC	12000	U	14000 U	U	NC	NC
1,1-Dichloroethane	10	μg/kg	270	U	100.00	U	NC	NC	3800	J	3700 J	J	2.67	NA
1,1-Dichloroethene	10	μg/kg	270	U	100.00	U	NC	NC	12000	U	14000 U	U	NC	NC
1,2,3-Trichlorobenzene	10	μg/kg	270	U	200.00	U	NC	NC	2600	J	27000 l	U	165	NA
1,2,4-Trichlorobenzene	10	μg/kg	140	J	200.00	U	35.3	NA	7000	J	11000 J	J	44.4	NA
1,2-Dibromo-3-chloropropane	10	μg/kg	270	U	300.00	U	NC	NC	12000	U	41000 U	U	NC	NC
1,2-Dibromoethane	10	μg/kg	270	U	100.00	U	NC	NC	12000	U	14000 U	U	NC	NC
1,2-Dichlorobenzene	10	μg/kg	2000		1600.00		22.2	NA	10000	J	17000 J	J	51.9	NA
1,2-Dichloroethane	10	μg/kg	270	U	100.00	U	NC	NC	12000	U	14000 l	U	NC	NC
1,2-Dichloropropane	10	μg/kg	270	U	100.00	U	NC	NC	12000	U	14000 l	U	NC	NC
1,3-Dichlorobenzene	10	μg/kg	610		560.00		8.55	NA	3900	J	5700 J	J	37.5	NA
1,4-Dichlorobenzene	10	μg/kg	2400		2300.00		4.26	NA	13000		22000 J	J	51.4	NA
1,4-Dioxane ²	10	μg/kg			10000	R	NC	NC						
2-Butanone	20	μg/kg	540	U	1000	U	NC	NC	25000	U	140000 l	U	NC	NC
2-Hexanone	20	μg/kg	540	U	1000	U	NC	NC	25000	U	140000 U	U	NC	NC
4-Methyl-2-pentanone	20	μg/kg	540	U	1000	U	NC	NC	25000	U	140000 U	U	NC	NC
Acetone	20	μg/kg	540	U	1000	U	NC	NC	25000	U	140000 U	U	NC	NC
Benzene	10	μg/kg	270	U	77		111	NA	12000	U	6800 L	U	NC	NC
Bromochloromethane	10	μg/kg	270	U	200.00	U	NC	NC	12000	U	27000 l	U	NC	NC
Bromodichloromethane	10	μg/kg	270	U	50.00	U	NC	NC	12000	U	6800 l	U	NC	NC
Bromoform	10	μg/kg	270	U	400.00	U	NC	NC	12000	U	54000 l	U	NC	NC
Bromomethane	10	μg/kg	270	U	200.00	U	NC	NC	12000	U	27000 l	U	NC	NC
Carbon Disulfide	10	μg/kg	270	U	1000	U	NC	NC	12000	U	140000 l	U	NC	NC
Carbon Tetrachloride	10	μg/kg	270	U	100.000	U	NC	NC	12000	U	14000 l	U	NC	NC
Chlorobenzene	10	μg/kg	16000		19000.000		17.1	NA	23000		36000		44.1	NA
Chloroethane	10	μg/kg	270	U	200.000	U	NC	NC	12000	U	27000 l	U	NC	NC
Chloroform	10	μg/kg	270	-	150.000	U	NC	NC	9600	J	20000 l	U	70.3	NA
Chloromethane	10	μg/kg	270	U	400.000	U	NC	NC	12000	U	54000 เ	U	NC	NC
cis-1,2-Dichloroethene	10	μg/kg	270	U	100.000	U	NC	NC	260000		230000		12.2	NA
cis-1,3-Dichloropropene	10	μg/kg	270	-	50.000	U	NC	NC	12000	U	6800 เ	U	NC	NC
Cyclohexane	10	μg/kg	270		1000		NC	NC	12000	U	140000 l	U	NC	NC
Dibromochloromethane	10	μg/kg	270	U	100.00	U	NC	NC	12000	U	14000 l	U	NC	NC



	2 x		GS-SS-111-003-CDM	GS-SS-111-003-G	EOSY			GS-B-114-001-CD	M	GS-B-114-001-GEO	SY		
Chemical Name		Unit	12/4/2018	12/4/2018		RPD	ABS	2/9/2019		2/9/2019		RPD	ABS
	CRQL		GS-SS	-111-003				GS-	-B-1	14-001			
Dichlorodifluoromethane	10	μg/kg	270 U	1000	U	NC	NC	12000	U	140000 l	U	NC	NC
Ethylbenzene	10	μg/kg	270 U	100	U	NC	NC	13000		19000		37.5	NA
Isopropylbenzene	10	μg/kg	270 U	100	U	NC	NC	12000	U	2300 J	J	136	NA
m,p-Xylene	10	μg/kg	270 U	200.00	J	NC	NC	29000		40000		31.9	NA
Methyl acetate	10	μg/kg	270 U	1500.00		139	NA	12000	U	54000 l	U	NC	NC
Methyl tert-Butyl Ether	10	μg/kg	270 U	200.00	J	NC	NC	12000	U	27000 l	U	NC	NC
Methylcyclohexane	10	μg/kg	270 U	400.00	J	NC	NC	12000	U	54000 l	U	NC	NC
Methylene Chloride	10	μg/kg	270 U	500.00	J	NC	NC	12000	U	68000 l	U	NC	NC
o-Xylene	10	μg/kg	270 U	100.00	J	NC	NC	9300	J	14000	•	40.3	NA
Styrene	10	μg/kg	270 U	100.00	J	NC	NC	12000	U	14000 l	U	NC	NC
Tetrachloroethene	10	μg/kg	270 U	50	J	NC	NC	1200000		2100000		54.5	NA
Toluene	10	μg/kg	270 U	100.00	J	NC	NC	3400	J	14000 l	U	122	NA
trans-1,2-Dichloroethene	10	μg/kg	270 U	150.00	U	NC	NC	12000	U	20000 l	U	NC	NC
trans-1,3-Dichloropropene	10	μg/kg	270 U	100.00	U	NC	NC	12000	U	14000 l	U	NC	NC
Trichloroethene	10	μg/kg	270 U	50	U	NC	NC	55000		66000		18.2	NA
Trichlorofluoromethane	10	μg/kg	270 U	400.00	U	NC	NC	12000	U	54000 l	U	NC	NC
Vinyl Chloride	10	μg/kg	270 U	100.00	U	NC	NC	5900	J	5600 J	J .	5.22	NA

Notes:

- 1. Sample pairs with RPD or ABS outside of criteria in the table are highlighted in yellow
- CDM Smith's assigned laboratory analyzed
 1,4-dioxane as a semivolatile organic compound.

ABS = absolute difference

CRQL = contract required quantitation limit

mg/kg = milligram per kilogram

NA = not applicable

NC = not calculable

RPD = relative percent difference

μg/kg = microgram per kilogram

VOC = volatile organic compound

Data Validation Qualifiers

J = The analyte was positively identified, and the associated numerical value is the approximate concentration of the analyte in the sample.

J+ = The result is an estimated quantity, but the result may be biased high.

J- = The result is an estimated quantity, but the result may be biased low.

NJ = The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents its approximate concentration.

R = The sample results are unusable due to poor data quality; certain criteria were not met. The analyte may or may not be present in the sample at the reported concentration.

U = The analyte was not detected at a level greater than or equal to the level of the sample CRQL.



Table 3-2
Groundwater Samples Data Comparison
Pierson's Creek Site, Operable Unit 2
Newark, New Jersey

	2 x		MW-10-CDI	M	MW-10-GEO	SY			MW-17-CDI	VI	MW-17-GEO	SY		
Chemical Name	CRQL	Unit	12/20/2018	8	12/19/2018	3	RPD	ABS	12/27/2018	8	12/27/2018	8	RPD	ABS
	CNQL		MW	/-10	-181220					ΜV	V-17			
Metals														
Aluminum	40	μg/L	39.4	J	10	J	NA	29.4	80.8		10	U	NA	70.8
Antimony	4	μg/L	2	U	4	J	NC	NC	2	U	4	J	NA	2
Arsenic	2	μg/L	68.7		20.77		107	NA	246		134.5		58.6	NA
Barium	20	μg/L	991		934.5		5.87	NA	750		782.4		4.23	NA
Beryllium	2	μg/L	1	U	0.5		NC	NC	1	U	0.5	U	NC	NC
Cadmium	2	μg/L	1	U	0.2	U	NC	NC	4.1		3.376		NA	0.72
Calcium	1000	μg/L	107000		105000		1.89	NA	111000		105000		5.56	NA
Chromium	4	μg/L	0.33	J	0.2051	J	NA	0.12	1		0.6716	J	NA	0.33
Cobalt	2	μg/L		U	0.3671	J	NA	0.63	950		829.6		13.5	NA
Copper	4	μg/L	2	U	1	J	NA	1	2.7		1.417		NA	1.28
Iron	400	μg/L	10500		592		NA	9908	9660		4430		74.2	NA
Lead	2	μg/L	0.5	U	1	U	NC	NC	22.2		4.667		NA	17.5
Magnesium	1000	μg/L	13900		12700		9.02	NA	18100		15300		16.8	NA
Manganese	2	μg/L	393		371.5		5.62	NA	1060		991.6		6.67	NA
Mercury	0.4	μg/L	0.05	UJ	0.2	J	NC	NC	54.2		6.195		159	NA
Nickel	2	μg/L	7		6.555		6.57	NA	8.5		6.798		22.3	NA
Potassium	1000	μg/L	20900		20100		3.9	NA	65500		59600		9.43	NA
Selenium	10	μg/L	3.5	J	5	J	NA	1.5	8.9		5	U	NA	3.9
Silver	2	μg/L	1	U	0.4	U	NC	NC	1	U	0.4	U	NC	NC
Sodium	1000	μg/L	250000		257000		2.76	NA	255000		239000		6.48	NA
Thallium	2	μg/L	1	U	0.5	J	NC	NC	1	U	0.2564	J	NA	0.74
Vanadium	10	μg/L	0.43	J	5	U	NA	4.57	1.4	J	5	U	NA	3.6
Zinc	4	μg/L	46.5		4.866	J	NA	41.6	880		577.5		41.5	NA
Polychlorinated Biphenyl														
Aroclor 1016	2	μg/L	1	U	0.25	U	NC	NC	1	U	0.25	U	NC	NC
Aroclor 1221	2	μg/L	1	U	0.25	J	NC	NC	1	U	0.25	U	NC	NC
Aroclor 1232	2	μg/L	1	U	0.25	U	NC	NC	1	U	0.25	U	NC	NC
Aroclor 1242	2	μg/L	1	U	0.25	U	NC	NC		U	0.25	U	NC	NC
Aroclor 1248	2	μg/L	1	U	0.25	U	NC	NC	1	U	0.25	U	NC	NC
Aroclor 1254	2	μg/L	1	U	0.25	U	NC	NC	1	U	0.25		NC	NC
Aroclor 1260	2	μg/L	1	U	0.25	U	NC	NC	1	U	0.25	U	NC	NC
Aroclor 1262	2	μg/L	1	U	0.25	U	NC	NC	1	U	0.25	U	NC	NC



Table 3-2
Groundwater Samples Data Comparison
Pierson's Creek Site, Operable Unit 2
Newark, New Jersey

	2 x		MW-10-CDM	MW-10-GEOSY			MW-17-CDM	MW-17-GEOSY		
Chemical Name	CROL	Unit	12/20/2018	12/19/2018	RPD	ABS	12/27/2018	12/27/2018	RPD	ABS
	CNQL		MW-10	-181220			MV	V-17		
Aroclor 1268	2	μg/L	1 U	0.25 U	NC	NC	1 U	0.25 U	NC	NC
Pesticides										
4,4'-DDD	0.2	μg/L	0.1 U	0.02 U	NC	NC	0.1 U	0.02 U	NC	NC
4,4'-DDE	0.2	μg/L	0.1 U	0.02 U	NC	NC	0.1 U	0.02 U	NC	NC
4,4'-DDT	0.2	μg/L	0.1 U	0.02 U	NC	NC	0.1 U	0.02 U	NC	NC
Aldrin	0.1	μg/L	0.05 U	0.01 U	NC	NC	0.05 U	0.01 U	NC	NC
alpha-BHC	0.1	μg/L	0.05 U	0.01 U	NC	NC	0.05 U	0.01 U	NC	NC
alpha-Chlordane	0.1	μg/L	0.05 U	0.01 U	NC	NC	0.05 U	0.01 U	NC	NC
beta-BHC	0.1	μg/L	0.05 U	0.01 U	NC	NC	0.05 U	0.01 U	NC	NC
delta-BHC	0.1	μg/L	0.05 U	0.01 U	NC	NC	0.05 U	0.21	NA	0.16
Dieldrin	0.1	μg/L	0.1 U	0.02 U	NC	NC	0.1 U	0.02 U	NC	NC
Endosulfan I	0.2	μg/L	0.05 U	0.01 U	NC	NC	0.05 U	0.01 U	NC	NC
Endosulfan II	0.1	μg/L	0.1 U	0.02 U	NC	NC	0.1 U	0.02 U	NC	NC
Endosulfan Sulfate	0.2	μg/L	0.1 U	0.02 U	NC	NC	0.1 U	0.02 U	NC	NC
Endrin	0.2	μg/L	0.1 U	0.02 U	NC	NC	0.1 U	0.02 U	NC	NC
Endrin aldehyde	0.2	μg/L	0.1 U	0.02 U	NC	NC	0.1 U	0.02 U	NC	NC
Endrin Ketone	0.2	μg/L	0.1 U	0.02 U	NC	NC	0.1 U	0.02 U	NC	NC
gamma-BHC (Lindane)	0.2	μg/L	0.05 UJ	0.01 U	NC	NC	0.05 U	0.01 U	NC	NC
gamma-Chlordane	0.1	μg/L	0.05 U	0.01 U	NC	NC	0.05 U	0.01 U	NC	NC
Heptachlor	0.1	μg/L	0.05 U	0.01 U	NC	NC	0.05 U	0.01 U	NC	NC
Heptachlor Epoxide	0.1	μg/L	0.05 U	0.01 U	NC	NC	0.05 U	0.01 U	NC	NC
Methoxychlor	1	μg/L	0.5 U	0.14 U	NC	NC	0.5 U	0.14 U	NC	NC
Toxaphene	10	μg/L	5 U	0.14 U	NC	NC	5 U	0.14 U	NC	NC
Semivolatile Organic Compound										
1,1'-Biphenyl	10	μg/L	5 U	2 UJ	NC	NC	5 U	2 U	NC	NC
1,2,4,5-Tetrachlorobenzene	10	μg/L	5 U	10 UJ	NC	NC	5 U	10 U	NC	NC
1,4-Dioxane ²	4	μg/L	2 U	250 R	NC	NC	2 UJ	500 R	NC	NC
2,2'-Oxybis(1-chloropropane)	20	μg/L	10 U	2 UJ	NC	NC	10 U	2 U	NC	NC
2,3,4,6-Tetrachlorophenol	10	μg/L	5 U	5 UJ	NC	NC	5 U	5 U	NC	NC
2,4,5-Trichlorophenol	10	μg/L	5 U	5 UJ	NC	NC	5 U	5 U	NC	NC
2,4,6-Trichlorophenol	10	μg/L	5 U	5 UJ	NC	NC	5 U	5 U	NC	NC
2,4-Dichlorophenol	10	μg/L	5 U	5 UJ	NC	NC	5 U	5 U	NC	NC



Table 3-2
Groundwater Samples Data Comparison
Pierson's Creek Site, Operable Unit 2
Newark, New Jersey

	2 x		MW-10-CDM	MW-10-GEO	SY			MW-17-CDM	MW-17-GEOS	Υ		
Chemical Name	CRQL	Unit	12/20/2018	12/19/2018	8	RPD	ABS	12/27/2018	12/27/2018		RPD	ABS
	CNQL		MW-10	0-181220					V-17			
2,4-Dimethylphenol	10	μg/L	5 U		UJ	NC	NC	5 U	5 L		NC	NC
2,4-Dinitrophenol	20	μg/L	10 U	20	U	NC	NC	10 U	20 L	J	NC	NC
2,4-Dinitrotoluene	10	μg/L	5 U	5	UJ	NC	NC	5 U	5 L	J	NC	NC
2,6-Dinitrotoluene	10	μg/L	5 U	5	UJ	NC	NC	5 U	5 L		NC	NC
2-Chloronaphthalene	10	μg/L	5 U	2	UJ	NC	NC	5 U	2 L	J	NC	NC
2-Chlorophenol	10	μg/L	5 U	2	UJ	NC	NC	5 U	2 L	J	NC	NC
2-Methylnaphthalene	10	μg/L	5 U		UJ	NC	NC	5 U	2 \		NC	NC
2-Methylphenol	20	μg/L	10 U	5	UJ	NC	NC	10 U	5 L		NC	NC
2-Nitroaniline	10	μg/L	5 U	5	UJ	NC	NC	5 U	5 L	JJ	NC	NC
2-Nitrophenol	10	μg/L	5 U	10	UJ	NC	NC	5 U	10 L	J	NC	NC
3,3'-Dichlorobenzidine	20	μg/L	10 U	5	UJ	NC	NC	10 U	5 L	J	NC	NC
3-Nitroaniline	20	μg/L	10 U	5	UJ	NC	NC	10 U	5 L	JJ	NC	NC
4,6-Dinitro-2-methylphenol	20	μg/L	10 U	0.69	UJ	NC	NC	10 U	0.69	JJ	NC	NC
4-Bromophenyl-phenylether	10	μg/L	5 U	2	UJ	NC	NC	5 U	2 L	J	NC	NC
4-Chloro-3-methylphenol	10	μg/L	5 U	2	UJ	NC	NC	5 U	2 L	J	NC	NC
4-Chloroaniline	20	μg/L	10 U	5	U	NC	NC	10 U	5 L	J	NC	NC
4-Chlorophenyl-phenylether	10	μg/L	5 U	2	UJ	NC	NC	5 U	2 \	J	NC	NC
4-Methylphenol	20	μg/L	10 U			NC	NC	10 U			NC	NC
4-Nitroaniline	20	μg/L	10 U	5	UJ	NC	NC	10 U	5 L	JJ	NC	NC
4-Nitrophenol	20	μg/L	10 U	10	U	NC	NC	10 U	10 L	J	NC	NC
Acenaphthene	10	μg/L	5 U	2	UJ	NC	NC	5 U	2 L	JJ	NC	NC
Acenaphthylene	10	μg/L	5 U	2	UJ	NC	NC	5 U	2 (J	NC	NC
Acetophenone	20	μg/L	10 U		UJ	NC	NC	10 U	5 L		NC	NC
Anthracene	10	μg/L	5 U	2	UJ	NC	NC	5 U	2 L	J	NC	NC
Atrazine	20	μg/L	10 U	3	UJ	NC	NC	10 U	3 L		NC	NC
Benzaldehyde	20	μg/L	10 U	5	U	NC	NC	10 U	5 L	J	NC	NC
Benzo(a)anthracene	10	μg/L	5 U	0.1	UJ	NC	NC	5 U	0.1	JJ	NC	NC
Benzo(a)pyrene	10	μg/L	5 U	0.1	UJ	NC	NC	5 U	0.1	JJ	NC	NC
Benzo(b)fluoranthene	10	μg/L	5 U	0.1	UJ	NC	NC	5 U	0.1	JJ	NC	NC
Benzo(g,h,i)perylene	10	μg/L	5 U	2	UJ	NC	NC	5 U	2 L	J	NC	NC
Benzo(k)fluoranthene	10	μg/L	5 U	0.1	UJ	NC	NC	5 U	0.1	JJ	NC	NC
Bis(2-chloroethoxy)methane	10	μg/L	5 U	5	UJ	NC	NC	5 U	5 L	J	NC	NC
Bis(2-chloroethyl)ether	20	μg/L	10 U	2	UJ	NC	NC	10 U	2 L	J	NC	NC



Table 3-2
Groundwater Samples Data Comparison
Pierson's Creek Site, Operable Unit 2
Newark, New Jersey

	2		MW-10-CDM	MW-10-GEOSY			MW-17-CDM	MW-17-GEOSY		
Chemical Name	2 x	Unit	12/20/2018	12/19/2018	RPD	ABS	12/27/2018	12/27/2018	RPD	ABS
	CRQL		MW-10	-181220			MV	V-17		
Bis(2-ethylhexyl)phthalate	10	μg/L	5 U	3 UJ	NC	NC	5 U	3 U	NC	NC
Butylbenzylphthalate	10	μg/L	5 U	5 UJ	NC	NC	5 U	5 UJ	NC	NC
Caprolactam	20	μg/L	10 U	10 U	NC	NC	10 U	10 U	NC	NC
Carbazole	20	μg/L	10 U	2 UJ	NC	NC	10 U	2 U	NC	NC
Chrysene	10	μg/L	5 U	2 UJ	NC	NC	5 U	2 U	NC	NC
Dibenzo(a,h)anthracene	10	μg/L	5 U	0.1 UJ	NC	NC	5 U	0.1 UJ	NC	NC
Dibenzofuran	10	μg/L	5 U	2 UJ	NC	NC	5 U	2 UJ	NC	NC
Diethylphthalate	10	μg/L	5 U	5 UJ	NC	NC	5 U	5 UJ	NC	NC
Dimethylphthalate	10	μg/L	5.5	5 UJ	NA	0.5	5 U	5 U	NC	NC
Di-n-butylphthalate	10	μg/L	5 U	5 UJ	NC	NC	5 U	5 UJ	NC	NC
Di-n-octylphthalate	20	μg/L	10 U	5 UJ	NC	NC	10 U	5 UJ	NC	NC
Fluoranthene	20	μg/L	10 U	2 UJ	NC	NC	10 U	2 U	NC	NC
Fluorene	10	μg/L	5 U	2 UJ	NC	NC	5 U	2 UJ	NC	NC
Hexachlorobenzene	10	μg/L	5 U	0.01 UJ	NC	NC	5 U	0.01 UJ	NC	NC
Hexachlorobutadiene	10	μg/L	5 U	1 UJ	NC	NC	5 U	1 U	NC	NC
Hexachlorocyclopentadiene	20	μg/L	10 U	20 U	NC	NC	10 U	20 U	NC	NC
Hexachloroethane	10	μg/L	5 U	2 U	NC	NC	5 U	2 U	NC	NC
Indeno(1,2,3-cd)pyrene	10	μg/L	5 U	0.1 UJ	NC	NC	5 U	0.1 UJ	NC	NC
Isophorone	10	μg/L	5 U	5 UJ	NC	NC	5 U	5 U	NC	NC
Naphthalene	10	μg/L	5 U	2 UJ	NC	NC	5 U	2 UJ	NC	NC
Nitrobenzene	10	μg/L	5 U	2 UJ	NC	NC	5 U	2 U	NC	NC
N-Nitroso-di-n-propylamine	10	μg/L	5 U	5 UJ	NC	NC	5 U	5 U	NC	NC
N-Nitrosodiphenylamine	10	μg/L	5 U	2 UJ	NC	NC	5 U	2 UJ	NC	NC
Pentachlorophenol	20	μg/L	10 U	0.3 J	NA	9.7	10 U	0.48	NA	9.52
Phenanthrene	10	μg/L	5 U	2 UJ	NC	NC	5 U	2 UJ	NC	NC
Phenol	20	μg/L	1.9 J	5 U	NA	3.1	10 U	0.61 J	NA	9.39
Pyrene	10	μg/L	5 U	2 UJ	NC	NC	5 U	2 UJ	NC	NC
Volatile Organic Compound										
1,1,1-Trichloroethane	1	μg/L	0.5 U	0.5 U	NC	NC	0.5 U	1 U	NC	NC
1,1,2,2-Tetrachloroethane	1	μg/L	0.5 U	0.5 U	NC	NC	0.5 U	1 U	NC	NC
1,1,2-Trichloro-1,2,2-trifluoroethane	1	μg/L	0.5 U	2.5 U	NC	NC	0.5 U	5 U	NC	NC
1,1,2-Trichloroethane	1	μg/L	0.5 U	0.75 U	NC	NC	0.5 U	1.5 U	NC	NC
1,1-Dichloroethane	1	μg/L	1.2	0.99	NA	0.21	1.4	1.3 J	NA	0.1



Table 3-2
Groundwater Samples Data Comparison
Pierson's Creek Site, Operable Unit 2
Newark, New Jersey

	2 x		MW-10-CDM	MW-10-GEOSY			MW-17-CDM	MW-17-GEOSY		
Chemical Name	CRQL	Unit	12/20/2018	12/19/2018	RPD	ABS	12/27/2018	12/27/2018	RPD	ABS
	CITQL		MW-10-181220				MW-17			
1,1-Dichloroethene	1	μg/L	0.5 U	0.5 U	NC	NC	0.5 U	1 U	NC	NC
1,2,3-Trichlorobenzene	1	μg/L	0.5 U	2.5 U	NC	NC	0.5 U	5 U	NC	NC
1,2,4-Trichlorobenzene	1	μg/L	0.5 U	2.5 U	NC	NC	0.5 U	5 U	NC	NC
1,2-Dibromo-3-chloropropane	1	μg/L	0.5 U	2.5 U	NC	NC	0.5 U	5 U	NC	NC
1,2-Dibromoethane	1	μg/L	0.5 U	2 U	NC	NC	0.5 U	4 U	NC	NC
1,2-Dichlorobenzene	1	μg/L	0.39 J	0.33 J	NA	0.06	0.35 J	5 U	NA	4.65
1,2-Dichloroethane	1	μg/L	0.5 U	0.5 U	NC	NC	0.5 U	0.37 J	NA	0.13
1,2-Dichloropropane	1	μg/L	0.5 U	1 U	NC	NC	0.5 U	2 U	NC	NC
1,3-Dichlorobenzene	1	μg/L	0.5 U	2.5 U	NC	NC	0.5 U	5 U	NC	NC
1,4-Dichlorobenzene	1	μg/L	0.5 U	2.5 U	NC	NC	0.5 U	5 U	NC	NC
1,4-Dioxane	4	μg/L		250 R	NC	NC		500 R	NC	NC
2-Butanone	10	μg/L	5 U	5 U	NC	NC	5 U	10 U	NC	NC
2-Hexanone	10	μg/L	5 U	5 U	NC	NC	5 U	10 U	NC	NC
4-Methyl-2-pentanone	10	μg/L	5 U	5 U	NC	NC	5 U	10 U	NC	NC
Acetone	10	μg/L	5 U	5 U	NC	NC	5 U	10 J	NA	5
Benzene	1	μg/L	6.5	5.8	11.4	NA	300	310	3.28	NA
Bromochloromethane	1	μg/L	0.5 U	2.5 U	NC	NC	0.5 U	5 U	NC	NC
Bromodichloromethane	1	μg/L	0.5 U	0.5 U	NC	NC	0.5 U	1 U	NC	NC
Bromoform	1	μg/L	0.5 U	2 U	NC	NC	0.5 U	4 U	NC	NC
Bromomethane	1	μg/L	0.5 U	1 U	NC	NC	0.5 U	2 U	NC	NC
Carbon Disulfide	1	μg/L	0.5 U	5 U	NC	NC	0.5 U	10 U	NC	NC
Carbon Tetrachloride	1	μg/L	0.5 U	0.5 U	NC	NC	0.5 U	1 U	NC	NC
Chlorobenzene	1	μg/L	0.29 J	0.34 J	NA	0.05	2.1	2.1	NA	0
Chloroethane	1	μg/L	0.77	0.77 J	NA	0	200	250	22.2	NA
Chloroform	1	μg/L	0.5 U	0.75 U	NC	NC	0.5 U	1.5 U	NC	NC
Chloromethane	1	μg/L	0.5 U	2.5 U	NC	NC	0.5 U	5 U	NC	NC
cis-1,2-Dichloroethene	1	μg/L	0.5 U	0.5 U	NC	NC	0.83	0.74 J	NA	0.09
cis-1,3-Dichloropropene	1	μg/L	0.5 U	0.5 U	NC	NC	0.5 U	1 U	NC	NC
Cyclohexane	1	μg/L	1.6	1.4 J	NA	0.2	0.42 J	20 U	NA	19.6
Dibromochloromethane	1	μg/L	0.5 U	0.5 U	NC	NC	0.5 U	1 U	NC	NC
Dichlorodifluoromethane	1	μg/L	0.5 U	5 U	NC	NC	0.5 U	10 U	NC	NC
Ethylbenzene	1	μg/L	0.33 J	0.28 J	NA	0.05	1	0.92 J	NA	0.08
Isopropylbenzene	1	μg/L	56	50	11.3	NA	8.5	6.3	29.7	NA



Table 3-2 Groundwater Samples Data Comparison Pierson's Creek Site, Operable Unit 2 Newark, New Jersey

	2 x CRQL	Unit	MW-10-CDN	VI	MW-10-GEOS	Υ			MW-17-CDN	VI	MW-17-GEOS	Υ		
Chemical Name			12/20/2018	/20/2018 12/19/2018			RPD	ABS	12/27/2018 12/27/2018			RPD	ABS	
			MW-10-181220					MW-17						
m,p-Xylene	1	μg/L	0.5	כ	1 U	J	NC	NC	0.56		2 l	J	NA	1.44
Methyl acetate	1	μg/L	0.5	J	2 U	J	NC	NC	0.5	J	4 l	J	NC	NC
Methyl tert-Butyl Ether	1	μg/L	0.5	כ	1 U	J	NC	NC	0.5	כ	2 l	J	NC	NC
Methylcyclohexane	1	μg/L	0.62		0.67 J		NA	0.05	0.21	J	20 L	J	NA	19.8
Methylene Chloride	1	μg/L	0.5	כ	2.5 U	J	NC	NC	0.54		5 L	J	NA	4.46
o-Xylene	1	μg/L	0.5	כ	1 U	J	NC	NC	1.6		1.2 J		NA	0.4
Styrene	1	μg/L	0.5	U	1 U	J	NC	NC	0.5	U	2 L	J	NC	NC
Tetrachloroethene	1	μg/L	0.5	כ	0.5 U	J	NC	NC	0.21	J	1 l	J	NA	0.79
Toluene	1	μg/L	0.14	J	0.75 U	J	NA	0.61	0.99		1.1 J		NA	0.11
trans-1,2-Dichloroethene	1	μg/L	0.5	כ	0.75 U	J	NC	NC	0.59		0.65 J		NA	0.06
trans-1,3-Dichloropropene	1	μg/L	0.5	כ	0.5 U	J	NC	NC	0.5	כ	1 l	J	NC	NC
Trichloroethene	1	μg/L	0.5	U	0.5 U	J	NC	NC	0.5	U	1 (J	NC	NC
Trichlorofluoromethane	1	μg/L	0.5	U	2.5 U	J	NC	NC	0.5	U	5 L	J	NC	NC
Vinyl Chloride	1	μg/L	0.5	U	0.2 U	J	NC	NC	0.5	U	0.5		NA	0

Notes:

- 1. No sample pairs are outside of RPD or ABS
- 2. PRP's laboratory analyzed 1,4-dioxane as a volatile organic compound.

ABS = absolute difference

CRQL = contract required quantitation limit

NA = not applicable

NC = not calculable

RPD = relative percent difference

μg/L = microgram per liter

Data Validation Qualifiers

J = The analyte was positively identified, and the associated numerical value is the approximate concentration of the analyte in the sample.

R = The sample results are unusable due to poor data quality; certain criteria were not met.

The analyte may or may not be present in the sample at the reported concentration.

U = The analyte was not detected at a level greater than or equal to the level of the sample CRQL.



Table 4-1 Soil Samples Statistical Data Comparison Summary Pierson's Creek Site, Operable Unit 2 Newark, New Jersey

Chemical	CAS No.	Number of Split Sample Pairs	Number of Split Sample Pairs with Detected Concentrations p-value (threshold 0.05)		Conclusion
Metals					
Antimony	7440-36-0	9	·		Data were not comparable with PRP sample results generally higher than CDM Smith split sample results
Arsenic	7440-38-2	9	9	0.063	Comparable
Silver	7440-22-4	9	5	0.032	Data were not comparable with PRP sample results generally higher than CDM Smith split sample results
Sodium	7440-23-5	9	8	0.419	Comparable
Thallium	7440-28-0	9	0	NC	All data were not detected
Vanadium	7440-62-2	9	9	0.399	Comparable
Zinc	7440-66-6	9	9	0.608	Comparable
Polychlorinated Biphenyl					
Aroclor 1242	53469-21-9	9	0	NC	All data were not detected
Aroclor 1254	11097-69-1	9	6	0.005	Data were not comparable with PRP sample results generally higher than CDM Smith split sample results
Aroclor 1260	11096-82-5	9	7	0.517	Comparable
Semivolatile Organic Compound					
1,1'-Biphenyl	92-52-4	9	0	NC	All data were not detected
1,2,4,5-Tetrachlorobenzene	95-94-3	9	0	NC	All data were not detected
2-Methylnaphthalene	91-57-6	9	2	NC	Less than 4 detected sample pairs
Acenaphthene	83-32-9	9	4	0.456	Comparable
Acenaphthylene	208-96-8	9	1	NC	Less than 4 detected sample pairs
Anthracene	120-12-7	9	5	0.124	Comparable
Benzo(a)anthracene	56-55-3	9	8	0.019	Data were not comparable with PRP sample results generally higher than CDM Smith split sample results



Table 4-1 Soil Samples Statistical Data Comparison Summary Pierson's Creek Site, Operable Unit 2 Newark, New Jersey

Chemical	CAS No.	Number of Split Sample Pairs	Number of Split Sample Pairs with Detected Concentrations	p-value (threshold 0.05)	Conclusion
Benzo(a)pyrene	50-32-8	9			Data were not comparable with PRP sample results generally higher than CDM Smith split sample results
Benzo(b)fluoranthene	205-99-2	9	9	0.024	Data were not comparable with PRP sample results generally higher than CDM Smith split sample results
Benzo(g,h,i)perylene	191-24-2	9	8	0.023	Data were not comparable with PRP sample results generally higher than CDM Smith split sample results
Benzo(k)fluoranthene	207-08-9	9	8	0.017	Data were not comparable with PRP sample results generally higher than CDM Smith split sample results
Bis(2-ethylhexyl)phthalate	117-81-7	9	5	0.193	Comparable
Carbazole	86-74-8	9	4	0.118	Comparable
Chrysene	218-01-9	9	8	0.024	Data were not comparable with PRP sample results generally higher than CDM Smith split sample results
Dibenzo(a,h)anthracene	53-70-3	9	4	0.108	Comparable
Dibenzofuran	132-64-9	9	3	NC	Less than 4 detected sample pairs
Diethylphthalate	84-66-2	9	0	NC	All data were not detected
Dimethylphthalate	131-11-3	9	0	NC	All data were not detected
Di-n-butylphthalate	84-74-2	9	2	NC	Less than 4 detected sample pairs
Di-n-octylphthalate	117-84-0	9	1	NC	Less than 4 detected sample pairs
Fluoranthene	206-44-0	9	7	0.020	Data were not comparable with PRP sample results generally higher than CDM Smith split sample results
Fluorene	86-73-7	9	5	0.318	Comparable
Indeno(1,2,3-cd)pyrene	193-39-5	9	7	0.009	Data were not comparable with PRP sample results generally higher than CDM Smith split sample results



Table 4-1 Soil Samples Statistical Data Comparison Summary Pierson's Creek Site, Operable Unit 2 Newark, New Jersey

Chemical	CAS No.	Number of Split Sample Pairs	Number of Split Sample Pairs with Detected Concentrations	p-value (threshold 0.05)	Conclusion
Naphthalene	91-20-3	9	4 0.094		Comparable
Phenanthrene	85-01-8	9	8	0.088	Comparable
Phenol	108-95-2	9	2	NC	Less than 4 detected sample pairs
Pyrene	129-00-0	9	8	0.056	Comparable
Volatile Organic Compound					
1,2,3-Trichlorobenzene	87-61-6	8	0	NC	All data were not detected
1,2,4-Trichlorobenzene	120-82-1	8	1	NC	Less than 4 detected sample pairs
1,4-Dichlorobenzene	106-46-7	8	3	NC	Less than 4 detected sample pairs
2-Butanone	78-93-3	8	0	NC	All data were not detected
Acetone	67-64-1	8	2	NC	Less than 4 detected sample pairs
Benzene	71-43-2	8	3	NC	Less than 4 detected sample pairs
Carbon Disulfide	75-15-0	8	0	NC	All data were not detected
Chlorobenzene	108-90-7	8	4	0.067	Comparable
cis-1,2-Dichloroethene	156-59-2	8	1	NC	Less than 4 detected sample pairs
Cyclohexane	110-82-7	8	2	NC	Less than 4 detected sample pairs
Ethylbenzene	100-41-4	8	5	0.164	Comparable
Isopropylbenzene	98-82-8	8	5	0.218	Comparable
m,p-Xylene	179601-23-1	8	4	0.069	Comparable
Methyl acetate	79-20-9	8	0	NC	All data were not detected
Methylcyclohexane	108-87-2	8	1	NC	Less than 4 detected sample pairs
o-Xylene	95-47-6	8	5	0.033	Data were not comparable with PRP sample results generally higher than CDM Smith split sample results
Tetrachloroethene	127-18-4	8	1	NC	Less than 4 detected sample pairs
Toluene	108-88-3	8	1	NC	Less than 4 detected sample pairs
Trichloroethene	79-01-6	8	1	NC	Less than 4 detected sample pairs

NC = not calculated for chemicals when less than four sample pairs were detected

PRP = potential resposible party

p-value from Paired Prentice Wilcoxon test less than 0.05 are **bolded**.

